

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP013123\
 Data File : BP013685.D
 Acq On : 31 Jan 2023 21:38
 Operator : CG/JU
 Sample : PB150501BS
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 PB150501BS

Manual Integrations
 APPROVED

Reviewed By : Christian Giraldo 02/01/2023
 Supervised By : Jagrut Upadhyay 02/01/2023

Quant Time: Feb 01 06:33:12 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP013123.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Feb 01 03:43:48 2023
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.140	152	266034	20.000	ng	0.00
21) Naphthalene-d8	10.969	136	927052	20.000	ng	0.00
39) Acenaphthene-d10	14.775	164	568324	20.000	ng	0.00
64) Phenanthrene-d10	17.528	188	1187851	20.000	ng	0.00
76) Chrysene-d12	21.604	240	1042325	20.000	ng	0.00
86) Perylene-d12	24.180	264	1120907	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.675	112	1669443	107.273	ng	0.00
7) Phenol-d6	7.287	99	2104936	101.525	ng	0.00
23) Nitrobenzene-d5	9.316	82	1142394	71.682	ng	0.00
42) 2,4,6-Tribromophenol	16.263	330	855169	100.322	ng	0.00
45) 2-Fluorobiphenyl	13.398	172	2965036	73.002	ng	0.00
79) Terphenyl-d14	20.057	244	3991337	76.040	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.516	88	290703	44.153	ng	99
3) Pyridine	3.934	79	744696	39.284	ng	99
4) n-Nitrosodimethylamine	3.834	42	294747	43.363	ng	99
6) Aniline	7.457	93	965938	35.737	ng	99
8) 2-Chlorophenol	7.699	128	750335	45.797	ng	99
9) Benzaldehyde	7.269	77	520438	55.455	ng	98
10) Phenol	7.316	94	979419	45.220	ng	98
11) bis(2-Chloroethyl)ether	7.552	93	760533	43.957	ng	99
12) 1,3-Dichlorobenzene	8.028	146	857690	46.473	ng	99
13) 1,4-Dichlorobenzene	8.175	146	881454	47.203	ng	99
14) 1,2-Dichlorobenzene	8.499	146	829972	46.448	ng	99
15) Benzyl Alcohol	8.381	79	631574	43.019	ng	99
16) 2,2'-oxybis(1-Chloropr...	8.675	45	878516	43.476	ng	100
17) 2-Methylphenol	8.581	107	624474	41.144	ng	99
18) Hexachloroethane	9.234	117	291416	47.326	ng	97
19) n-Nitroso-di-n-propyla...	8.951	70	532311	40.921	ng	99
20) 3+4-Methylphenols	8.910	107	836830	40.416	ng	98
22) Acetophenone	8.969	105	1080877	44.711	ng	99
24) Nitrobenzene	9.357	77	784485	45.382	ng	99
25) Isophorone	9.887	82	1467206	41.536	ng	99
26) 2-Nitrophenol	10.075	139	345936	43.489	ng	99
27) 2,4-Dimethylphenol	10.122	122	665002	46.649	ng	99
28) bis(2-Chloroethoxy)met...	10.363	93	934499	44.139	ng	99
29) 2,4-Dichlorophenol	10.610	162	721224	46.482	ng	100
30) 1,2,4-Trichlorobenzene	10.828	180	819887	47.622	ng	99
31) Naphthalene	11.022	128	2162258	44.359	ng	99
32) Benzoic acid	10.251	122	435698	39.012	ng	100
33) 4-Chloroaniline	11.122	127	317124	14.806	ng	98
34) Hexachlorobutadiene	11.304	225	521517	47.721	ng	98
35) Caprolactam	11.910	113	192523	37.782	ng	98
36) 4-Chloro-3-methylphenol	12.234	107	673356	41.256	ng	100
37) 2-Methylnaphthalene	12.616	142	1524752	41.602	ng	99
38) 1-Methylnaphthalene	12.834	142	1403146	40.755	ng	98
40) 1,2,4,5-Tetrachloroben...	12.975	216	946907	48.803	ng	99
41) Hexachlorocyclopentadiene	12.957	237	1178883	133.119	ng	98
43) 2,4,6-Trichlorophenol	13.210	196	600467	46.998	ng	100

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP013123\
 Data File : BP013685.D
 Acq On : 31 Jan 2023 21:38
 Operator : CG/JU
 Sample : PB150501BS
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 PB150501BS

Manual Integrations
 APPROVED

Reviewed By : Christian Giraldo 02/01/2023
 Supervised By : Jagrut Upadhyay 02/01/2023

Quant Time: Feb 01 06:33:12 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP013123.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Feb 01 03:43:48 2023
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	13.281	196	646426	43.213	ng	99
46) 1,1'-Biphenyl	13.610	154	1992609	46.332	ng	99
47) 2-Chloronaphthalene	13.657	162	1577340	47.448	ng	99
48) 2-Nitroaniline	13.857	65	322437	39.497	ng	98
49) Acenaphthylene	14.498	152	2385413	44.496	ng	100
50) Dimethylphthalate	14.228	163	1852807	42.149	ng	99
51) 2,6-Dinitrotoluene	14.345	165	371428	41.551	ng	97
52) Acenaphthene	14.839	154	1660969m	48.125	ng	
53) 3-Nitroaniline	14.675	138	219819	24.317	ng	93
54) 2,4-Dinitrophenol	14.881	184	438023	74.937	ng	99
55) Dibenzofuran	15.175	168	2324721	43.039	ng	99
56) 4-Nitrophenol	14.975	139	643626	84.448	ng	99
57) 2,4-Dinitrotoluene	15.128	165	486942	40.842	ng	96
58) Fluorene	15.822	166	1803001	42.515	ng	99
59) 2,3,4,6-Tetrachlorophenol	15.392	232	572572	44.018	ng	99
60) Diethylphthalate	15.586	149	1702279	40.395	ng	99
61) 4-Chlorophenyl-phenylether	15.810	204	986706	42.726	ng	99
62) 4-Nitroaniline	15.839	138	343562	36.464	ng	99
63) Azobenzene	16.104	77	1639972	42.128	ng	99
65) 4,6-Dinitro-2-methylph...	15.892	198	272660	40.606	ng	97
66) n-Nitrosodiphenylamine	16.028	169	1565837	45.305	ng	99
67) 4-Bromophenyl-phenylether	16.710	248	628136	46.130	ng	98
68) Hexachlorobenzene	16.828	284	735010	48.747	ng	99
69) Atrazine	16.975	200	553721	49.235	ng	100
70) Pentachlorophenol	17.169	266	910661	81.877	ng	100
71) Phenanthrene	17.569	178	2809058	44.120	ng	99
72) Anthracene	17.663	178	2867210	44.606	ng	99
73) Carbazole	17.927	167	2497697	42.793	ng	99
74) Di-n-butylphthalate	18.457	149	2661317	41.667	ng	100
75) Fluoranthene	19.522	202	3215198	40.563	ng	99
77) Benzidine	19.692	184	1517406	67.399	ng	99
78) Pyrene	19.874	202	3347276	47.733	ng	100
80) Butylbenzylphthalate	20.727	149	862762	45.067	ng	99
81) Benzo(a)anthracene	21.586	228	3208130	44.312	ng	100
82) 3,3'-Dichlorobenzidine	21.510	252	787247	34.394	ng	98
83) Chrysene	21.645	228	3032152	43.959	ng	100
84) Bis(2-ethylhexyl)phtha...	21.486	149	1386742	42.544	ng	99
85) Di-n-octyl phthalate	22.468	149	2458328	41.301	ng	100
87) Indeno(1,2,3-cd)pyrene	26.898	276	3900672	45.758	ng	100
88) Benzo(b)fluoranthene	23.386	252	3176495	47.151	ng	100
89) Benzo(k)fluoranthene	23.439	252	3239692	48.011	ng	100
90) Benzo(a)pyrene	24.068	252	2841860	42.497	ng	99
91) Dibenzo(a,h)anthracene	26.921	278	3248567	45.910	ng	99
92) Benzo(g,h,i)perylene	27.739	276	3196971	45.490	ng	100

(#) = qualifier out of range (m) = manual integration (+) = signals summed

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP013123\
 Data File : BP013685.D
 Acq On : 31 Jan 2023 21:38
 Operator : CG/JU
 Sample : PB150501BS
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :

BNA_P

ClientSampleId :

PB150501BS

Manual Integrations

APPROVED

Reviewed By : Christian Giraldo 02/01/2023
 Supervised By : Jagrut Upadhyay 02/01/2023

Quant Time: Feb 01 06:33:12 2023

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP013123.M

Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

QLast Update : Wed Feb 01 03:43:48 2023

Response via : Initial Calibration

