

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP012723\
 Data File : BP013612.D
 Acq On : 27 Jan 2023 11:45
 Operator : CG/JU
 Sample : PB150456BSD
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 PB150456BSD

Manual Integrations
 APPROVED

Reviewed By : Christian Giraldo 01/30/2023
 Supervised By : Jagrut Upadhyay 01/30/2023

Quant Time: Jan 29 23:48:30 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP012623.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jan 27 03:27:54 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.158	152	200490	20.000	ng	0.00
21) Naphthalene-d8	10.987	136	706794	20.000	ng	0.00
39) Acenaphthene-d10	14.793	164	376335	20.000	ng	0.00
64) Phenanthrene-d10	17.545	188	705429	20.000	ng	0.00
76) Chrysene-d12	21.627	240	619854	20.000	ng	0.00
86) Perylene-d12	24.216	264	730195	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.687	112	1615076	144.354	ng	0.00
7) Phenol-d6	7.299	99	1943859	128.613	ng	0.00
23) Nitrobenzene-d5	9.328	82	1129994	87.848	ng	0.00
42) 2,4,6-Tribromophenol	16.281	330	637388	111.244	ng	0.00
45) 2-Fluorobiphenyl	13.416	172	2553972	96.925	ng	0.00
79) Terphenyl-d14	20.075	244	3111769	88.588	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.528	88	200877	40.722	ng	98
3) Pyridine	3.940	79	522040	37.230	ng	99
4) n-Nitrosodimethylamine	3.846	42	236554	43.166	ng	97
6) Aniline	7.469	93	690413	34.468	ng	99
8) 2-Chlorophenol	7.711	128	572326	48.188	ng	98
9) Benzaldehyde	7.281	77	275794m	27.322	ng	
10) Phenol	7.328	94	705756	44.578	ng	99
11) bis(2-Chloroethyl)ether	7.563	93	588800	44.909	ng	99
12) 1,3-Dichlorobenzene	8.046	146	665436	48.537	ng	98
13) 1,4-Dichlorobenzene	8.193	146	672644	48.384	ng	99
14) 1,2-Dichlorobenzene	8.511	146	644229	49.220	ng	99
15) Benzyl Alcohol	8.393	79	467142m	39.804	ng	
16) 2,2'-oxybis(1-Chloropr...	8.681	45	669991	45.348	ng	99
17) 2-Methylphenol	8.593	107	467872	40.954	ng	98
18) Hexachloroethane	9.252	117	236360	47.457	ng	98
19) n-Nitroso-di-n-propyla...	8.969	70	386834	39.384	ng	98
20) 3+4-Methylphenols	8.922	107	607290	38.790	ng	98
22) Acetophenone	8.987	105	810871	45.451	ng	# 98
24) Nitrobenzene	9.369	77	594650	44.172	ng	100
25) Isophorone	9.899	82	1114092	39.540	ng	100
26) 2-Nitrophenol	10.087	139	249588	43.159	ng	96
27) 2,4-Dimethylphenol	10.140	122	468917	43.803	ng	98
28) bis(2-Chloroethoxy)met...	10.381	93	698360	43.562	ng	99
29) 2,4-Dichlorophenol	10.622	162	519650	43.136	ng	100
30) 1,2,4-Trichlorobenzene	10.846	180	633431	47.425	ng	98
31) Naphthalene	11.034	128	1604887	44.685	ng	100
32) Benzoic acid	10.257	122	271586	34.801	ng	92
33) 4-Chloroaniline	11.140	127	250365	16.057	ng	98
34) Hexachlorobutadiene	11.322	225	418153	46.563	ng	98
35) Caprolactam	11.916	113	110995	32.133	ng	97
36) 4-Chloro-3-methylphenol	12.246	107	465708	36.466	ng	98
37) 2-Methylnaphthalene	12.628	142	1078089	39.277	ng	98
38) 1-Methylnaphthalene	12.846	142	992045	38.615	ng	98
40) 1,2,4,5-Tetrachloroben...	12.993	216	675673	51.535	ng	99
41) Hexachlorocyclopentadiene	12.975	237	817570	120.695	ng	98
43) 2,4,6-Trichlorophenol	13.222	196	405485	46.221	ng	99

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	13.293	196	418161	40.759	ng	98
46) 1,1'-Biphenyl	13.628	154	1369518	48.835	ng	99
47) 2-Chloronaphthalene	13.669	162	1063800	49.707	ng	99
48) 2-Nitroaniline	13.869	65	159113	32.379	ng	96
49) Acenaphthylene	14.516	152	1578964	44.730	ng	99
50) Dimethylphthalate	14.240	163	1164046	39.821	ng	99
51) 2,6-Dinitrotoluene	14.357	165	210543	36.044	ng	97
52) Acenaphthene	14.857	154	1059667m	46.903	ng	
53) 3-Nitroaniline	14.687	138	129043	23.787	ng	96
54) 2,4-Dinitrophenol	14.887	184	247646	68.945	ng	97
55) Dibenzofuran	15.187	168	1492296	42.174	ng	98
56) 4-Nitrophenol	14.981	139	372392	78.152	ng	93
57) 2,4-Dinitrotoluene	15.140	165	270973	34.685	ng	98
58) Fluorene	15.840	166	1108169	40.954	ng	99
59) 2,3,4,6-Tetrachlorophenol	15.410	232	351225	39.518	ng	99
60) Diethylphthalate	15.598	149	1038641	38.907	ng	99
61) 4-Chlorophenyl-phenylether	15.828	204	615305	40.270	ng	97
62) 4-Nitroaniline	15.851	138	175660	33.653	ng	92
63) Azobenzene	16.122	77	1062767	39.072	ng	98
65) 4,6-Dinitro-2-methylph...	15.904	198	141856	37.399	ng	97
66) n-Nitrosodiphenylamine	16.040	169	946064	48.676	ng	97
67) 4-Bromophenyl-phenylether	16.728	248	380366	46.780	ng	98
68) Hexachlorobenzene	16.845	284	440644	49.579	ng	97
69) Atrazine	16.992	200	289369	48.705	ng	99
70) Pentachlorophenol	17.187	266	545143	78.925	ng	99
71) Phenanthrene	17.586	178	1710219	47.270	ng	99
72) Anthracene	17.681	178	1704453	46.624	ng	100
73) Carbazole	17.939	167	1485913	45.614	ng	99
74) Di-n-butylphthalate	18.481	149	1284669	43.327	ng	99
75) Fluoranthene	19.539	202	1927098	42.705	ng	99
77) Benzidine	19.710	184	319641	49.185	ng	98
78) Pyrene	19.892	202	2024646	42.205	ng	99
80) Butylbenzylphthalate	20.745	149	227374	35.981	ng	97
81) Benzo(a)anthracene	21.610	228	1975559	45.260	ng	100
82) 3,3'-Dichlorobenzidine	21.533	252	380831	36.355	ng	98
83) Chrysene	21.669	228	1851795	45.151	ng	99
84) Bis(2-ethylhexyl)phtha...	21.516	149	379922	39.142	ng	100
85) Di-n-octyl phthalate	22.504	149	726218	45.302	ng	98
87) Indeno(1,2,3-cd)pyrene	26.957	276	2954999	57.563	ng	99
88) Benzo(b)fluoranthene	23.421	252	2173180	46.357	ng	99
89) Benzo(k)fluoranthene	23.474	252	2096879	45.057	ng	99
90) Benzo(a)pyrene	24.104	252	1940884	43.987	ng	99
91) Dibenzo(a,h)anthracene	26.974	278	2464687	57.700	ng	99
92) Benzo(g,h,i)perylene	27.792	276	2488137	59.158	ng	99

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

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