

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM111222\
 Data File : BM037491.D
 Acq On : 12 Nov 2022 14:03
 Operator : CG/JU
 Sample : SSTDICC020
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTDICC020

Manual Integrations
 APPROVED

Reviewed By : Christian Giraldo 11/14/2022
 Supervised By : Jagrut Upadhyay 11/14/2022

Quant Time: Nov 14 00:36:15 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM111222.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Nov 14 00:33:11 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.798	152	168306	20.000	ng	0.00
21) Naphthalene-d8	10.598	136	783135	20.000	ng	0.00
39) Acenaphthene-d10	14.445	164	525581	20.000	ng	0.00
64) Phenanthrene-d10	17.192	188	1114903	20.000	ng	0.00
76) Chrysene-d12	21.380	240	1303844	20.000	ng	0.00
86) Perylene-d12	23.715	264	1398427	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.369	112	360354	36.990	ng	0.00
7) Phenol-d6	6.981	99	528530	39.974	ng	0.00
23) Nitrobenzene-d5	8.969	82	594608	38.307	ng	0.00
42) 2,4,6-Tribromophenol	15.939	330	282996	45.690	ng	0.00
45) 2-Fluorobiphenyl	13.069	172	1461464	36.947	ng	0.00
79) Terphenyl-d14	19.821	244	2599486	35.753	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.246	88	75495	18.925	ng	99
3) Pyridine	3.663	79	237094	19.950	ng	98
4) n-Nitrosodimethylamine	3.575	42	107124	20.622	ng	# 96
6) Aniline	7.134	93	326640	20.208	ng	99
8) 2-Chlorophenol	7.363	128	239246	20.178	ng	99
9) Benzaldehyde	6.951	77	150130m	22.432	ng	
10) Phenol	7.004	94	294478	19.877	ng	98
11) bis(2-Chloroethyl)ether	7.234	93	238127	20.047	ng	99
12) 1,3-Dichlorobenzene	7.687	146	256615	19.806	ng	98
13) 1,4-Dichlorobenzene	7.834	146	265432	19.861	ng	98
14) 1,2-Dichlorobenzene	8.151	146	264026	20.569	ng	99
15) Benzyl Alcohol	8.057	79	197123	20.853	ng	98
16) 2,2'-oxybis(1-Chloropr...	8.328	45	354586	20.939	ng	98
17) 2-Methylphenol	8.257	107	208960	21.582	ng	98
18) Hexachloroethane	8.869	117	99498	21.110	ng	98
19) n-Nitroso-di-n-propyla...	8.616	70	206651	22.723	ng	99
20) 3+4-Methylphenols	8.586	107	297908	22.401	ng	98
22) Acetophenone	8.634	105	396076	19.458	ng	# 99
24) Nitrobenzene	9.016	77	303747	19.302	ng	99
25) Isophorone	9.539	82	541695	20.881	ng	98
26) 2-Nitrophenol	9.722	139	138575	19.670	ng	99
27) 2,4-Dimethylphenol	9.786	122	199789	19.109	ng	99
28) bis(2-Chloroethoxy)met...	10.028	93	314798	19.503	ng	100
29) 2,4-Dichlorophenol	10.257	162	241420	20.115	ng	99
30) 1,2,4-Trichlorobenzene	10.463	180	262250	19.268	ng	99
31) Naphthalene	10.651	128	868164	19.567	ng	100
32) Benzoic acid	9.916	122	159468	18.799	ng	100
33) 4-Chloroaniline	10.780	127	349477	19.774	ng	99
34) Hexachlorobutadiene	10.922	225	151908	19.988	ng	98
35) Caprolactam	11.575	113	84367	23.122	ng	95
36) 4-Chloro-3-methylphenol	11.910	107	263505	21.305	ng	98
37) 2-Methylnaphthalene	12.263	142	603706	20.082	ng	99
38) 1-Methylnaphthalene	12.486	142	598312	20.369	ng	99
40) 1,2,4,5-Tetrachloroben...	12.633	216	290064	18.355	ng	100
41) Hexachlorocyclopentadiene	12.604	237	108542	14.416	ng	99
43) 2,4,6-Trichlorophenol	12.886	196	209935	19.298	ng	99

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM111222\
 Data File : BM037491.D
 Acq On : 12 Nov 2022 14:03
 Operator : CG/JU
 Sample : SSTDICC020
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTDICC020

Manual Integrations
 APPROVED

Reviewed By : Christian Giraldo 11/14/2022
 Supervised By : Jagrut Upadhyay 11/14/2022

Quant Time : Nov 14 00:36:15 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM111222.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Nov 14 00:33:11 2022
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	12.957	196	231282	19.273	ng	99
46) 1,1'-Biphenyl	13.280	154	792444	18.684	ng	98
47) 2-Chloronaphthalene	13.321	162	640542	18.978	ng	99
48) 2-Nitroaniline	13.545	65	194080	19.898	ng	98
49) Acenaphthylene	14.168	152	1022506	19.516	ng	100
50) Dimethylphthalate	13.916	163	810790	20.197	ng	99
51) 2,6-Dinitrotoluene	14.039	165	175614	20.657	ng	97
52) Acenaphthene	14.510	154	617633	19.068	ng	98
53) 3-Nitroaniline	14.374	138	186770	19.665	ng	97
54) 2,4-Dinitrophenol	14.586	184	87949	18.599	ng	96
55) Dibenzofuran	14.845	168	998550	20.015	ng	100
56) 4-Nitrophenol	14.692	139	138328	18.257	ng	98
57) 2,4-Dinitrotoluene	14.827	165	239010	21.365	ng	97
58) Fluorene	15.492	166	836017	20.076	ng	100
59) 2,3,4,6-Tetrachlorophenol	15.074	232	213200	21.196	ng	98
60) Diethylphthalate	15.274	149	833295	21.357	ng	100
61) 4-Chlorophenyl-phenylether	15.492	204	387219	20.098	ng	98
62) 4-Nitroaniline	15.533	138	172897	18.151	ng	98
63) Azobenzene	15.786	77	800289	20.820	ng	100
65) 4,6-Dinitro-2-methylph...	15.592	198	135858	18.966	ng	99
66) n-Nitrosodiphenylamine	15.710	169	697276	18.503	ng	99
67) 4-Bromophenyl-phenylether	16.386	248	239623	19.152	ng	100
68) Hexachlorobenzene	16.492	284	295958	19.547	ng	99
69) Atrazine	16.668	200	193631	20.421	ng	97
70) Pentachlorophenol	16.845	266	172613	19.592	ng	97
71) Phenanthrene	17.233	178	1310595	18.947	ng	100
72) Anthracene	17.327	178	1276610	19.495	ng	100
73) Carbazole	17.604	167	1192689	19.856	ng	100
74) Di-n-butylphthalate	18.162	149	1492583	22.504	ng	100
75) Fluoranthene	19.251	202	1531793	20.950	ng	99
77) Benzidine	19.456	184	270400m	13.908	ng	
78) Pyrene	19.615	202	1645612	16.413	ng	99
80) Butylbenzylphthalate	20.515	149	698309	19.503	ng	98
81) Benzo(a)anthracene	21.362	228	1775382	19.126	ng	99
82) 3,3'-Dichlorobenzidine	21.303	252	570783	20.936	ng	100
83) Chrysene	21.415	228	1720068	19.233	ng	98
84) Bis(2-ethylhexyl)phtha...	21.292	149	1028954	21.411	ng	98
85) Di-n-octyl phthalate	22.192	149	1781461	23.538	ng	99
87) Indeno(1,2,3-cd)pyrene	26.132	276	2221410	19.401	ng	99
88) Benzo(b)fluoranthene	23.003	252	1767830	18.209	ng	99
89) Benzo(k)fluoranthene	23.050	252	1818359	18.362	ng	99
90) Benzo(a)pyrene	23.609	252	1486297	18.789	ng	99
91) Dibenzo(a,h)anthracene	26.144	278	1826108	19.205	ng	100
92) Benzo(g,h,i)perylene	26.868	276	1817742	19.642	ng	99

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

