

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM080822\
 Data File : BM036146.D
 Acq On : 08 Aug 2022 14:56
 Operator : CG/JU
 Sample : N4063-03MSD
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 P-5CMSD

Manual Integrations
 APPROVED

Reviewed By : Christian Giraldo 08/09/2022
 Supervised By : Jagrut Upadhyay 08/09/2022

Quant Time: Aug 08 23:33:17 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM080122.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Aug 05 03:36:25 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.851	152	371721	20.000	ng	0.00
21) Naphthalene-d8	10.657	136	1593799	20.000	ng	0.00
39) Acenaphthene-d10	14.498	164	968309	20.000	ng	0.00
64) Phenanthrene-d10	17.239	188	1907076	20.000	ng	0.00
76) Chrysene-d12	21.427	240	1466661	20.000	ng	0.00
86) Perylene-d12	23.792	264	1207550	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.440	112	2635623	114.953	ng	0.02
7) Phenol-d6	7.069	99	3377535	115.772	ng	0.04
23) Nitrobenzene-d5	9.028	82	2010907	70.626	ng	0.00
42) 2,4,6-Tribromophenol	15.992	330	1292711	113.800	ng	0.00
45) 2-Fluorobiphenyl	13.122	172	4608145	70.265	ng	0.00
79) Terphenyl-d14	19.868	244	6459765	86.874	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.275	88	337066	36.483	ng	91
3) Pyridine	3.687	79	955121	38.552	ng	89
4) n-Nitrosodimethylamine	3.605	42	439194	43.140	ng	# 64
6) Aniline	7.187	93	1544681	47.813	ng	95
8) 2-Chlorophenol	7.428	128	1196805	48.957	ng	95
9) Benzaldehyde	6.993	77	752639	48.730	ng	95
10) Phenol	7.098	94	1457843	49.628	ng	90
11) bis(2-Chloroethyl)ether	7.281	93	1087415	44.937	ng	96
12) 1,3-Dichlorobenzene	7.734	146	1229630	45.392	ng	97
13) 1,4-Dichlorobenzene	7.887	146	1263282	45.728	ng	98
14) 1,2-Dichlorobenzene	8.204	146	1209572	45.378	ng	99
15) Benzyl Alcohol	8.116	79	978066m	49.959	ng	
16) 2,2'-oxybis(1-Chloropr...	8.393	45	1467031	45.592	ng	96
17) 2-Methylphenol	8.340	107	977885	50.289	ng	95
18) Hexachloroethane	8.928	117	471431	47.405	ng	97
19) n-Nitroso-di-n-propyla...	8.675	70	817407	46.127	ng	87
20) 3+4-Methylphenols	8.669	107	1292346	48.918	ng	# 87
22) Acetophenone	8.687	105	1698063	45.008	ng	# 94
24) Nitrobenzene	9.069	77	1228319	45.890	ng	95
25) Isophorone	9.598	82	2333487	50.379	ng	98
26) 2-Nitrophenol	9.775	139	643037	50.934	ng	92
27) 2,4-Dimethylphenol	9.863	122	1092163	55.845	ng	98
28) bis(2-Chloroethoxy)met...	10.081	93	1457276	44.160	ng	99
29) 2,4-Dichlorophenol	10.328	162	1088777	49.505	ng	100
30) 1,2,4-Trichlorobenzene	10.522	180	1076577	43.195	ng	98
31) Naphthalene	10.710	128	3549450	44.794	ng	99
32) Benzoic acid	10.063	122	822940	53.084	ng	93
33) 4-Chloroaniline	10.834	127	1228184	38.459	ng	96
34) Hexachlorobutadiene	10.986	225	648403	44.273	ng	97
35) Caprolactam	11.651	113	348179	51.397	ng	93
36) 4-Chloro-3-methylphenol	11.992	107	1144382	49.504	ng	99
37) 2-Methylnaphthalene	12.322	142	2437944	46.234	ng	98
38) 1-Methylnaphthalene	12.539	142	2339686	45.309	ng	98
40) 1,2,4,5-Tetrachloroben...	12.686	216	1190823	45.361	ng	98
41) Hexachlorocyclopentadiene	12.663	237	1489693	107.143	ng	99
43) 2,4,6-Trichlorophenol	12.945	196	843557	47.223	ng	99

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	13.028	196	912056	48.367	ng	97
46) 1,1'-Biphenyl	13.333	154	3171537	44.943	ng	100
47) 2-Chloronaphthalene	13.375	162	2427362	44.943	ng	99
48) 2-Nitroaniline	13.598	65	728325	49.483	ng	89
49) Acenaphthylene	14.222	152	3865482	46.300	ng	100
50) Dimethylphthalate	13.969	163	2958840	44.934	ng	99
51) 2,6-Dinitrotoluene	14.092	165	652918	49.419	ng	90
52) Acenaphthene	14.563	154	2738438m	50.142	ng	
53) 3-Nitroaniline	14.427	138	604803	39.236	ng	93
54) 2,4-Dinitrophenol	14.627	184	707354	101.174	ng	91
55) Dibenzofuran	14.898	168	3600099	44.020	ng	98
56) 4-Nitrophenol	14.780	139	1224581	99.274	ng	87
57) 2,4-Dinitrotoluene	14.874	165	894279	51.641	ng	98
58) Fluorene	15.545	166	2910334	45.113	ng	98
59) 2,3,4,6-Tetrachlorophenol	15.133	232	740331	46.347	ng	97
60) Diethylphthalate	15.327	149	3068204	45.341	ng	99
61) 4-Chlorophenyl-phenyle...	15.539	204	1412052	43.896	ng	98
62) 4-Nitroaniline	15.592	138	751130m	47.367	ng	
63) Azobenzene	15.833	77	2852944	44.348	ng	94
65) 4,6-Dinitro-2-methylph...	15.633	198	449724	46.120	ng	94
66) n-Nitrosodiphenylamine	15.763	169	2488191	46.218	ng	99
67) 4-Bromophenyl-phenylether	16.433	248	867480	45.476	ng	96
68) Hexachlorobenzene	16.539	284	1015582	46.401	ng	94
69) Atrazine	16.721	200	880295	58.437	ng	98
70) Pentachlorophenol	16.898	266	1195050	93.228	ng	99
71) Phenanthrene	17.286	178	4416005	45.451	ng	99
72) Anthracene	17.374	178	4446942	47.276	ng	98
73) Carbazole	17.657	167	4150143	43.545	ng	99
74) Di-n-butylphthalate	18.215	149	5357327	47.088	ng	99
75) Fluoranthene	19.304	202	4891972	44.981	ng	99
77) Benzidine	19.498	184	3379557	153.647	ng	99
78) Pyrene	19.662	202	4940214	54.319	ng	98
80) Butylbenzylphthalate	20.568	149	2255996	57.473	ng	95
81) Benzo(a)anthracene	21.415	228	4215148	46.660	ng	99
82) 3,3'-Dichlorobenzidine	21.351	252	1371860	62.541	ng	99
83) Chrysene	21.468	228	3884500	45.235	ng	100
84) Bis(2-ethylhexyl)phtha...	21.339	149	3155036	53.277	ng	99
85) Di-n-octyl phthalate	22.256	149	4659041	48.207	ng	# 95
87) Indeno(1,2,3-cd)pyrene	26.250	276	3772048	44.450	ng	99
88) Benzo(b)fluoranthene	23.074	252	3588322	50.719	ng	99
89) Benzo(k)fluoranthene	23.121	252	3352217	46.912	ng	99
90) Benzo(a)pyrene	23.686	252	2985824	50.300	ng	99
91) Dibenzo(a,h)anthracene	26.268	278	3297561	46.423	ng	98
92) Benzo(g,h,i)perylene	27.009	276	3281029	47.713	ng	99

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

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