

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM102623\
 Data File : BM042455.D
 Acq On : 26 Oct 2023 16:36
 Operator : MA/JU
 Sample : PB156601BS
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 PB156601BS

Quant Time: Oct 26 17:17:25 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM102123.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Oct 21 02:32:53 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.998	152	134691	20.000	ng	0.00
21) Naphthalene-d8	10.822	136	531083	20.000	ng	0.00
39) Acenaphthene-d10	14.645	164	297714	20.000	ng	0.00
64) Phenanthrene-d10	17.392	188	586030	20.000	ng	0.00
76) Chrysene-d12	21.574	240	444274	20.000	ng	0.00
86) Perylene-d12	24.038	264	466755	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.528	112	1215378	127.530	ng	0.00
7) Phenol-d6	7.157	99	1546312	121.075	ng	0.00
23) Nitrobenzene-d5	9.198	82	1016371	84.626	ng	0.00
42) 2,4,6-Tribromophenol	16.139	330	541806	162.316	ng	0.00
45) 2-Fluorobiphenyl	13.269	172	2026268	91.652	ng	0.00
79) Terphenyl-d14	20.003	244	2779539	110.956	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.393	88	134029	29.867	ng	96
3) Pyridine	3.816	79	361647	27.468	ng	95
4) n-Nitrosodimethylamine	3.740	42	229102	34.464	ng	91
6) Aniline	7.339	93	481615	29.618	ng	97
8) 2-Chlorophenol	7.557	128	421257	44.547	ng	94
9) Benzaldehyde	7.157	77	73759	9.916	ng	99
10) Phenol	7.181	94	539878	40.910	ng	98
11) bis(2-Chloroethyl)ether	7.434	93	430799	39.761	ng	97
12) 1,3-Dichlorobenzene	7.881	146	445970	45.595	ng	96
13) 1,4-Dichlorobenzene	8.039	146	456858	46.401	ng	97
14) 1,2-Dichlorobenzene	8.345	146	436721	46.193	ng	98
15) Benzyl Alcohol	8.257	79	393495	41.563	ng	97
16) 2,2'-oxybis(1-Chloropr...	8.528	45	676127	39.438	ng	99
17) 2-Methylphenol	8.439	107	354375	40.693	ng	98
18) Hexachloroethane	9.063	117	180638	47.657	ng	# 90
19) n-Nitroso-di-n-propyla...	8.816	70	350627	40.506	ng	98
20) 3+4-Methylphenols	8.775	107	471009	40.324	ng	98
22) Acetophenone	8.851	105	633734	43.740	ng	# 95
24) Nitrobenzene	9.245	77	512294	42.636	ng	97
25) Isophorone	9.751	82	929285	45.530	ng	98
26) 2-Nitrophenol	9.945	139	227270	49.080	ng	94
27) 2,4-Dimethylphenol	9.980	122	356551	43.067	ng	97
28) bis(2-Chloroethoxy)met...	10.239	93	566636	43.819	ng	98
29) 2,4-Dichlorophenol	10.463	162	388005	53.892	ng	96
30) 1,2,4-Trichlorobenzene	10.675	180	414452	53.697	ng	97
31) Naphthalene	10.875	128	1271577	46.614	ng	100
32) Benzoic acid	10.128	122	298771	49.192	ng	98
33) 4-Chloroaniline	11.010	127	291078	24.577	ng	98
34) Hexachlorobutadiene	11.098	225	254781	61.002	ng	99
35) Caprolactam	11.839	113	122053	43.361	ng	89
36) 4-Chloro-3-methylphenol	12.104	107	412814	47.514	ng	97
37) 2-Methylnaphthalene	12.474	142	838873	45.683	ng	98
38) 1-Methylnaphthalene	12.692	142	813891	47.311	ng	99
40) 1,2,4,5-Tetrachloroben...	12.827	216	453957	52.744	ng	97
41) Hexachlorocyclopentadiene	12.769	237	431957	91.844	ng	99
43) 2,4,6-Trichlorophenol	13.080	196	299559	53.766	ng	97

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	13.151	196	324929	49.334	ng	# 93
46) 1,1'-Biphenyl	13.480	154	1082851	45.849	ng	98
47) 2-Chloronaphthalene	13.533	162	826665	47.709	ng	98
48) 2-Nitroaniline	13.763	65	291151	38.788	ng	91
49) Acenaphthylene	14.374	152	1363206	48.732	ng	100
50) Dimethylphthalate	14.110	163	1012946	46.386	ng	99
51) 2,6-Dinitrotoluene	14.251	165	220356	44.890	ng	85
52) Acenaphthene	14.710	154	777765	45.988	ng	99
53) 3-Nitroaniline	14.592	138	181214	32.151	ng	96
54) 2,4-Dinitrophenol	14.798	184	287020	92.547	ng	90
55) Dibenzofuran	15.045	168	1238334	46.294	ng	98
56) 4-Nitrophenol	14.880	139	397872	89.458	ng	93
57) 2,4-Dinitrotoluene	15.033	165	289801	44.288	ng	86
58) Fluorene	15.692	166	987415	46.779	ng	98
59) 2,3,4,6-Tetrachlorophenol	15.257	232	278760	56.340	ng	89
60) Diethylphthalate	15.451	149	992442	46.053	ng	97
61) 4-Chlorophenyl-phenyle...	15.680	204	506013	50.904	ng	93
62) 4-Nitroaniline	15.751	138	213886	38.906	ng	97
63) Azobenzene	15.974	77	1089471	42.138	ng	95
65) 4,6-Dinitro-2-methylph...	15.786	198	166466	44.918	ng	# 79
66) n-Nitrosodiphenylamine	15.904	169	849076	48.103	ng	98
67) 4-Bromophenyl-phenylether	16.574	248	303249	55.832	ng	95
68) Hexachlorobenzene	16.662	284	348251	59.984	ng	# 91
69) Atrazine	16.851	200	295643	66.048	ng	99
70) Pentachlorophenol	17.021	266	465521	94.848	ng	99
71) Phenanthrene	17.439	178	1499425	47.033	ng	100
72) Anthracene	17.527	178	1506989	47.729	ng	99
73) Carbazole	17.815	167	1369594	46.239	ng	99
74) Di-n-butylphthalate	18.333	149	1729876	46.231	ng	99
75) Fluoranthene	19.445	202	1697712	48.353	ng	98
77) Benzidine	19.656	184	537888	83.704	ng	99
78) Pyrene	19.815	202	1743824	56.170	ng	100
80) Butylbenzylphthalate	20.692	149	741708	52.386	ng	89
81) Benzo(a)anthracene	21.556	228	1471418	49.404	ng	99
82) 3,3'-Dichlorobenzidine	21.497	252	505661	56.874	ng	# 97
83) Chrysene	21.615	228	1341702	46.903	ng	100
84) Bis(2-ethylhexyl)phtha...	21.433	149	1032185	51.341	ng	97
85) Di-n-octyl phthalate	22.380	149	1711267	49.486	ng	95
87) Indeno(1,2,3-cd)pyrene	26.632	276	1676389	49.146	ng	94
88) Benzo(b)fluoranthene	23.274	252	1407038	50.081	ng	98
89) Benzo(k)fluoranthene	23.327	252	1284785	44.046	ng	98
90) Benzo(a)pyrene	23.927	252	1213183	44.863	ng	98
91) Dibenzo(a,h)anthracene	26.650	278	1395024	49.190	ng	97
92) Benzo(g,h,i)perylene	27.444	276	1376850	49.991	ng	# 93

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

