

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM042121\
 Data File : BM029570.D
 Acq On : 21 Apr 2021 15:03
 Operator : CG/JU
 Sample : SSTDICV040
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 ICVBM042121

Quant Time: Apr 21 15:54:32 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM042121.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Apr 21 15:14:18 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.645	152	71822	20.00	ng	0.00
21) Naphthalene-d8	10.428	136	287683	20.00	ng	0.00
39) Acenaphthene-d10	14.292	164	215631	20.00	ng	0.00
64) Phenanthrene-d10	17.033	188	478916	20.00	ng	0.00
76) Chrysene-d12	21.221	240	564409	20.00	ng	0.00
86) Perylene-d12	23.415	264	585166	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.246	112	290896	81.20	ng	0.00
7) Phenol-d6	6.828	99	444796	83.13	ng	0.00
23) Nitrobenzene-d5	8.804	82	469188	84.00	ng	0.00
42) 2,4,6-Tribromophenol	15.786	330	252951	84.82	ng	0.00
45) 2-Fluorobiphenyl	12.910	172	1314735	82.75	ng	0.00
79) Terphenyl-d14	19.668	244	2388460	80.23	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.175	88	56635	39.21	ng	99
3) Pyridine	3.581	79	155866	42.34	ng	98
4) n-Nitrosodimethylamine	3.493	42	70262	38.90	ng	98
6) Aniline	6.981	93	246731	41.94	ng	99
8) 2-Chlorophenol	7.216	128	181198	40.68	ng	99
9) Benzaldehyde	6.798	77	118857	39.07	ng	98
10) Phenol	6.851	94	213994	41.27	ng	100
11) bis(2-Chloroethyl)ether	7.081	93	163203	41.26	ng	99
12) 1,3-Dichlorobenzene	7.539	146	206783	40.28	ng	98
13) 1,4-Dichlorobenzene	7.681	146	210348	40.02	ng	99
14) 1,2-Dichlorobenzene	7.998	146	210258	40.13	ng	96
15) Benzyl Alcohol	7.892	79	165025	42.45	ng	98
16) 2,2'-oxybis(1-Chloropr...	8.181	45	207240	39.74	ng	99
17) 2-Methylphenol	8.092	107	151282	41.87	ng	98
18) Hexachloroethane	8.716	117	76868	40.50	ng	100
19) n-Nitroso-di-n-propyla...	8.457	70	163191	41.82	ng	96
20) 3+4-Methylphenols	8.422	107	210048	42.26	ng	98
22) Acetophenone	8.469	105	290077	42.06	ng	# 98
24) Nitrobenzene	8.845	77	235089	41.62	ng	98
25) Isophorone	9.369	82	438969	41.82	ng	100
26) 2-Nitrophenol	9.551	139	107759	40.23	ng	99
27) 2,4-Dimethylphenol	9.616	122	159754	42.37	ng	97
28) bis(2-Chloroethoxy)met...	9.851	93	216251	41.70	ng	100
29) 2,4-Dichlorophenol	10.080	162	211982	42.62	ng	97
30) 1,2,4-Trichlorobenzene	10.292	180	235350	41.09	ng	99
31) Naphthalene	10.480	128	599376	40.52	ng	99
32) Benzoic acid	9.769	122	132652	41.35	ng	99
33) 4-Chloroaniline	10.592	127	250777	42.99	ng	98
34) Hexachlorobutadiene	10.763	225	150246	41.20	ng	95
35) Caprolactam	11.386	113	64614	42.90	ng	95
36) 4-Chloro-3-methylphenol	11.722	107	207304	42.65	ng	99
37) 2-Methylnaphthalene	12.098	142	483856	40.86	ng	100
38) 1-Methylnaphthalene	12.322	142	461365	40.86	ng	99
40) 1,2,4,5-Tetrachloroben...	12.469	216	272089	42.08	ng	99
41) Hexachlorocyclopentadiene	12.451	237	149672	41.48	ng	100
43) 2,4,6-Trichlorophenol	12.716	196	193482	43.52	ng	100

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM042121\
 Data File : BM029570.D
 Acq On : 21 Apr 2021 15:03
 Operator : CG/JU
 Sample : SSTDICV040
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 ICVBM042121

Quant Time: Apr 21 15:54:32 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM042121.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Apr 21 15:14:18 2021
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	12.786	196	210664	42.92	ng	97
46) 1,1'-Biphenyl	13.121	154	654337	41.51	ng	98
47) 2-Chloronaphthalene	13.163	162	525908	41.69	ng	99
48) 2-Nitroaniline	13.369	65	142185	40.79	ng	95
49) Acenaphthylene	14.010	152	812814	42.03	ng	99
50) Dimethylphthalate	13.757	163	684843	42.02	ng	100
51) 2,6-Dinitrotoluene	13.874	165	154270	43.00	ng	96
52) Acenaphthene	14.357	154	502510	41.48	ng	99
53) 3-Nitroaniline	14.204	138	135822	41.23	ng	98
54) 2,4-Dinitrophenol	14.410	184	89330	41.30	ng	90
55) Dibenzofuran	14.692	168	848806	41.49	ng	97
56) 4-Nitrophenol	14.515	139	115438	43.50	ng	96
57) 2,4-Dinitrotoluene	14.663	165	214347	44.24	ng	98
58) Fluorene	15.345	166	719961	41.92	ng	100
59) 2,3,4,6-Tetrachlorophenol	14.921	232	190554	43.64	ng	98
60) Diethylphthalate	15.127	149	672324	42.03	ng	98
61) 4-Chlorophenyl-phenyle...	15.339	204	360510	41.48	ng	97
62) 4-Nitroaniline	15.368	138	139834	39.75	ng	98
63) Azobenzene	15.633	77	643497	42.37	ng	99
65) 4,6-Dinitro-2-methylph...	15.427	198	134029	42.10	ng	98
66) n-Nitrosodiphenylamine	15.557	169	612609	41.38	ng	99
67) 4-Bromophenyl-phenylether	16.233	248	210453	42.08	ng	98
68) Hexachlorobenzene	16.345	284	233617	41.09	ng	99
69) Atrazine	16.509	200	228885	43.33	ng	98
70) Pentachlorophenol	16.686	266	156261	42.21	ng	96
71) Phenanthrene	17.080	178	1129160	40.95	ng	99
72) Anthracene	17.168	178	1131163	41.58	ng	100
73) Carbazole	17.439	167	1068002	41.90	ng	99
74) Di-n-butylphthalate	18.009	149	1175322	42.47	ng	100
75) Fluoranthene	19.092	202	1388100	41.51	ng	100
77) Benzidine	19.286	184	532935	44.60	ng	100
78) Pyrene	19.456	202	1464395	40.33	ng	100
80) Butylbenzylphthalate	20.368	149	568189	41.80	ng	95
81) Benzo(a)anthracene	21.203	228	1494136	40.01	ng	99
82) 3,3'-Dichlorobenzidine	21.144	252	486887	40.82	ng	98
83) Chrysene	21.256	228	1460096	39.79	ng	100
84) Bis(2-ethylhexyl)phtha...	21.144	149	907575	41.66	ng	99
85) Di-n-octyl phthalate	22.009	149	1514496	41.43	ng	100
87) Indeno(1,2,3-cd)pyrene	25.627	276	1867201	41.27	ng	100
88) Benzo(b)fluoranthene	22.762	252	1554218	40.73	ng	99
89) Benzo(k)fluoranthene	22.803	252	1550555	40.54	ng	99
90) Benzo(a)pyrene	23.321	252	1447294	40.40	ng	100
91) Dibenzo(a,h)anthracene	25.638	278	1614674	41.39	ng	100
92) Benzo(g,h,i)perylene	26.303	276	1491836	41.35	ng	99

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

