

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM091720\
 Data File : BM027455.D
 Acq On : 17 Sep 2020 19:54
 Operator : JU/CG
 Sample : SSTDICV040
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 ICVBM091720

Quant Time: Sep 17 20:27:04 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\8270-BM091720.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Sep 17 20:07:45 2020
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.504	152	172310	20.00	ng	0.00
21) Naphthalene-d8	10.269	136	613336	20.00	ng	0.00
39) Acenaphthene-d10	14.151	164	405989	20.00	ng	0.00
64) Phenanthrene-d10	16.898	188	880739	20.00	ng	0.00
76) Chrysene-d12	21.103	240	911137	20.00	ng	0.00
86) Perylene-d12	23.244	264	883825	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.122	112	742935	77.27	ng	0.00
7) Phenol-d6	6.698	99	1015332	78.41	ng	0.00
23) Nitrobenzene-d5	8.651	82	1092248	75.71	ng	0.00
42) 2,4,6-Tribromophenol	15.645	330	497257	75.18	ng	0.00
45) 2-Fluorobiphenyl	12.769	172	2340808	76.03	ng	0.00
79) Terphenyl-d14	19.551	244	3938294	77.12	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.087	88	176496	35.80	ng	99
3) Pyridine	3.469	79	464567	35.64	ng	98
4) n-Nitrosodimethylamine	3.387	42	173838	35.91	ng	# 94
6) Aniline	6.840	93	570896	38.85	ng	99
8) 2-Chlorophenol	7.075	128	379943	38.32	ng	99
9) Benzaldehyde	6.651	77	337504	38.83	ng	97
10) Phenol	6.722	94	481266	38.88	ng	96
11) bis(2-Chloroethyl)ether	6.940	93	410599	38.42	ng	99
12) 1,3-Dichlorobenzene	7.392	146	491047	37.76	ng	99
13) 1,4-Dichlorobenzene	7.534	146	497404	37.92	ng	98
14) 1,2-Dichlorobenzene	7.851	146	472626	37.75	ng	98
15) Benzyl Alcohol	7.745	79	408876	39.00	ng	99
16) 2,2'-oxybis(1-Chloropr...	8.045	45	317502	38.28	ng	99
17) 2-Methylphenol	7.951	107	320876	38.86	ng	99
18) Hexachloroethane	8.569	117	201515	37.75	ng	97
19) n-Nitroso-di-n-propyla...	8.316	70	370253	37.49	ng	97
20) 3+4-Methylphenols	8.281	107	430999	39.34	ng	98
22) Acetophenone	8.316	105	615415	37.24	ng	98
24) Nitrobenzene	8.692	77	552666	37.64	ng	99
25) Isophorone	9.216	82	856745	37.51	ng	98
26) 2-Nitrophenol	9.392	139	211700	39.97	ng	96
27) 2,4-Dimethylphenol	9.469	122	285761	38.59	ng	95
28) bis(2-Chloroethoxy)met...	9.704	93	555803	38.66	ng	97
29) 2,4-Dichlorophenol	9.928	162	395293	38.43	ng	99
30) 1,2,4-Trichlorobenzene	10.139	180	506269	37.70	ng	99
31) Naphthalene	10.322	128	1185700	37.45	ng	99
32) Benzoic acid	9.628	122	218246	38.88	ng	97
33) 4-Chloroaniline	10.433	127	482841	38.84	ng	98
34) Hexachlorobutadiene	10.616	225	355687	36.70	ng	99
35) Caprolactam	11.222	113	102820	39.58	ng	97
36) 4-Chloro-3-methylphenol	11.575	107	396003	37.99	ng	98
37) 2-Methylnaphthalene	11.939	142	894434	37.20	ng	99
38) 1-Methylnaphthalene	12.163	142	847745	36.98	ng	100
40) 1,2,4,5-Tetrachloroben...	12.322	216	564998	38.60	ng	99
41) Hexachlorocyclopentadiene	12.310	237	316299	40.37	ng	97
43) 2,4,6-Trichlorophenol	12.569	196	355666	39.19	ng	95

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44) 2,4,5-Trichlorophenol	12.639	196	363969	39.38	ng	97
46) 1,1'-Biphenyl	12.980	154	1188400	37.99	ng	99
47) 2-Chloronaphthalene	13.016	162	997312	37.98	ng	98
48) 2-Nitroaniline	13.222	65	310841	39.99	ng	100
49) Acenaphthylene	13.869	152	1429871	38.34	ng	100
50) Dimethylphthalate	13.621	163	1226869	36.99	ng	99
51) 2,6-Dinitrotoluene	13.733	165	260396	38.58	ng	97
52) Acenaphthene	14.216	154	887892	37.45	ng	98
53) 3-Nitroaniline	14.063	138	245903	40.27	ng	97
54) 2,4-Dinitrophenol	14.269	184	144220	39.07	ng	97
55) Dibenzofuran	14.551	168	1451748	36.80	ng	100
56) 4-Nitrophenol	14.380	139	187435	39.64	ng	97
57) 2,4-Dinitrotoluene	14.527	165	359085	38.90	ng	97
58) Fluorene	15.204	166	1193085	36.73	ng	99
59) 2,3,4,6-Tetrachlorophenol	14.786	232	350797	37.60	ng	99
60) Diethylphthalate	14.998	149	1236971	36.80	ng	99
61) 4-Chlorophenyl-phenyle...	15.210	204	687225	36.43	ng	98
62) 4-Nitroaniline	15.227	138	258281	38.35	ng	97
63) Azobenzene	15.498	77	1250181	36.79	ng	97
65) 4,6-Dinitro-2-methylph...	15.298	198	189867	39.31	ng	97
66) n-Nitrosodiphenylamine	15.421	169	1012055	38.61	ng	98
67) 4-Bromophenyl-phenylether	16.098	248	438106	38.56	ng	97
68) Hexachlorobenzene	16.215	284	506694	38.07	ng	97
69) Atrazine	16.386	200	378886	38.84	ng	99
70) Pentachlorophenol	16.557	266	275140	39.16	ng	98
71) Phenanthrene	16.939	178	1887080	37.58	ng	100
72) Anthracene	17.033	178	1846388	38.57	ng	99
73) Carbazole	17.304	167	1620417	39.25	ng	99
74) Di-n-butylphthalate	17.892	149	2111214	39.04	ng	99
75) Fluoranthene	18.968	202	2249749	38.14	ng	99
77) Benzidine	19.156	184	773755	39.65	ng	99
78) Pyrene	19.327	202	2315167	38.93	ng	100
80) Butylbenzylphthalate	20.256	149	932949	39.99	ng	97
81) Benzo(a)anthracene	21.086	228	2274628	38.42	ng	99
82) 3,3'-Dichlorobenzidine	21.027	252	806828	40.91	ng	99
83) Chrysene	21.139	228	2192814	37.80	ng	99
84) Bis(2-ethylhexyl)phtha...	21.050	149	1398152	39.63	ng	99
85) Di-n-octyl phthalate	21.903	149	2257886	39.31	ng	100
87) Indeno(1,2,3-cd)pyrene	25.374	276	2565021	38.30	ng	99
88) Benzo(b)fluoranthene	22.615	252	2275413	38.27	ng	100
89) Benzo(k)fluoranthene	22.656	252	2264426	38.63	ng	100
90) Benzo(a)pyrene	23.156	252	2091304	38.94	ng	100
91) Dibenzo(a,h)anthracene	25.380	278	2120653	38.23	ng	99
92) Benzo(g,h,i)perylene	26.015	276	2055054	38.15	ng	99

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

