

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM082521\
 Data File : BM031891.D
 Acq On : 25 Aug 2021 14:14
 Operator : CG/JU
 Sample : SSTDICV040
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 ICVBM082521

Quant Time: Aug 25 15:26:22 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM082521.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Aug 25 14:16:52 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.522	152	38500	20.000	ng	0.00
21) Naphthalene-d8	10.281	136	174467	20.000	ng	0.00
39) Acenaphthene-d10	14.163	164	123441	20.000	ng	0.00
64) Phenanthrene-d10	16.916	188	268565	20.000	ng	0.00
76) Chrysene-d12	21.127	240	230975	20.000	ng	0.00
86) Perylene-d12	23.268	264	179083	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.157	112	178011	76.439	ng	0.00
7) Phenol-d6	6.722	99	254948	77.374	ng	0.00
23) Nitrobenzene-d5	8.675	82	299187	74.280	ng	0.00
42) 2,4,6-Tribromophenol	15.669	330	113204	77.582	ng	0.00
45) 2-Fluorobiphenyl	12.775	172	687053	73.432	ng	0.00
79) Terphenyl-d14	19.568	244	1160522	72.528	ng	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.169	88	33202	39.724	ng	# 85
3) Pyridine	3.552	79	88925	37.517	ng	# 85
4) n-Nitrosodimethylamine	3.469	42	63503	39.440	ng	# 46
6) Aniline	6.869	93	134798	38.512	ng	97
8) 2-Chlorophenol	7.093	128	101968	37.609	ng	96
9) Benzaldehyde	6.687	77	78033	42.601	ng	97
10) Phenol	6.746	94	124417	38.182	ng	97
11) bis(2-Chloroethyl)ether	6.969	93	91519	38.502	ng	93
12) 1,3-Dichlorobenzene	7.410	146	113048	37.752	ng	97
13) 1,4-Dichlorobenzene	7.557	146	114495	37.316	ng	97
14) 1,2-Dichlorobenzene	7.863	146	111758	37.408	ng	97
15) Benzyl Alcohol	7.769	79	109134	38.826	ng	96
16) 2,2'-oxybis(1-Chloropr...	8.040	45	119697	37.535	ng	95
17) 2-Methylphenol	7.969	107	88492	38.065	ng	98
18) Hexachloroethane	8.569	117	47864	37.658	ng	95
19) n-Nitroso-di-n-propyla...	8.334	70	98573	38.830	ng	97
20) 3+4-Methylphenols	8.298	107	123634	38.656	ng	98
22) Acetophenone	8.346	105	175208	36.944	ng	97
24) Nitrobenzene	8.716	77	149692	36.705	ng	99
25) Isophorone	9.240	82	258910	37.221	ng	98
26) 2-Nitrophenol	9.410	139	63449	38.863	ng	90
27) 2,4-Dimethylphenol	9.481	122	91090	37.072	ng	96
28) bis(2-Chloroethoxy)met...	9.716	93	126872	37.341	ng	99
29) 2,4-Dichlorophenol	9.940	162	109157	37.827	ng	99
30) 1,2,4-Trichlorobenzene	10.145	180	118880	37.549	ng	98
31) Naphthalene	10.328	128	359388	36.862	ng	99
32) Benzoic acid	9.675	122	81285	38.829	ng	93
33) 4-Chloroaniline	10.457	127	147635	37.878	ng	98
34) Hexachlorobutadiene	10.604	225	76114	36.349	ng	97
35) Caprolactam	11.281	113	34641	39.391	ng	# 83
36) 4-Chloro-3-methylphenol	11.592	107	131321	38.375	ng	99
37) 2-Methylnaphthalene	11.951	142	276088	38.034	ng	98
38) 1-Methylnaphthalene	12.175	142	255433	37.202	ng	99
40) 1,2,4,5-Tetrachloroben...	12.328	216	134076	37.075	ng	98
41) Hexachlorocyclopentadiene	12.292	237	66093	37.483	ng	99
43) 2,4,6-Trichlorophenol	12.581	196	97907	37.705	ng	96

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44) 2,4,5-Trichlorophenol	12.651	196	105705	37.670	ng	97
46) 1,1'-Biphenyl	12.986	154	351636	36.807	ng	99
47) 2-Chloronaphthalene	13.022	162	275659	36.649	ng	99
48) 2-Nitroaniline	13.251	65	99343	37.955	ng	94
49) Acenaphthylene	13.880	152	447385	36.975	ng	100
50) Dimethylphthalate	13.639	163	362128	37.007	ng	100
51) 2,6-Dinitrotoluene	13.763	165	82526	38.295	ng	88
52) Acenaphthene	14.227	154	269494	36.584	ng	99
53) 3-Nitroaniline	14.092	138	80760	38.527	ng	100
54) 2,4-Dinitrophenol	14.310	184	50686	41.530	ng	88
55) Dibenzofuran	14.563	168	460378	37.112	ng	99
56) 4-Nitrophenol	14.416	139	68206	41.182	ng	94
57) 2,4-Dinitrotoluene	14.557	165	120489	39.256	ng	93
58) Fluorene	15.222	166	382320	37.585	ng	100
59) 2,3,4,6-Tetrachlorophenol	14.798	232	94003	38.372	ng	96
60) Diethylphthalate	15.016	149	397064	37.179	ng	98
61) 4-Chlorophenyl-phenyle...	15.222	204	188102	37.253	ng	97
62) 4-Nitroaniline	15.269	138	82541	40.775	ng	96
63) Azobenzene	15.516	77	440313	37.391	ng	99
65) 4,6-Dinitro-2-methylph...	15.327	198	73147	39.543	ng	88
66) n-Nitrosodiphenylamine	15.445	169	321473	35.914	ng	97
67) 4-Bromophenyl-phenylether	16.116	248	107668	35.893	ng	96
68) Hexachlorobenzene	16.221	284	114763	36.197	ng	94
69) Atrazine	16.410	200	114419	37.919	ng	98
70) Pentachlorophenol	16.574	266	76902	38.619	ng	98
71) Phenanthrene	16.963	178	579377	36.641	ng	100
72) Anthracene	17.051	178	571181	36.931	ng	99
73) Carbazole	17.333	167	548311	37.540	ng	99
74) Di-n-butylphthalate	17.904	149	711450	38.802	ng	100
75) Fluoranthene	18.986	202	670933	38.853	ng	98
77) Benzidine	19.192	184	250054	41.386	ng	99
78) Pyrene	19.351	202	682389	35.326	ng	99
80) Butylbenzylphthalate	20.274	149	315009	37.918	ng	99
81) Benzo(a)anthracene	21.109	228	619046	37.261	ng	100
82) 3,3'-Dichlorobenzidine	21.056	252	185211	38.770	ng	99
83) Chrysene	21.162	228	596151	37.218	ng	100
84) Bis(2-ethylhexyl)phtha...	21.056	149	478190	40.008	ng	98
85) Di-n-octyl phthalate	21.915	149	709148	39.372	ng	97
87) Indeno(1,2,3-cd)pyrene	25.386	276	439247	35.447	ng	97
88) Benzo(b)fluoranthene	22.639	252	529126	39.025	ng	97
89) Benzo(k)fluoranthene	22.680	252	498753	38.805	ng	98
90) Benzo(a)pyrene	23.180	252	444487	37.849	ng	99
91) Dibenzo(a,h)anthracene	25.397	278	379959	35.403	ng	98
92) Benzo(g,h,i)perylene	26.033	276	349314	34.421	ng	100

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

