

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM012121\
 Data File : BM028489.D
 Acq On : 21 Jan 2021 15:16
 Operator : CG/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTDCCC040EC

Quant Time: Jan 21 15:50:33 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM012021.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jan 20 14:45:01 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.028	152	41794	20.00	ng	0.00
21) Naphthalene-d8	10.845	136	179416	20.00	ng	# 0.00
39) Acenaphthene-d10	14.675	164	104185	20.00	ng	0.00
64) Phenanthrene-d10	17.398	188	187542	20.00	ng	# 0.00
76) Chrysene-d12	21.527	240	134920	20.00	ng	# 0.00
86) Perylene-d12	23.815	264	125921	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.552	112	186794	70.60	ng	0.00
7) Phenol-d6	7.193	99	274541	76.21	ng	0.00
23) Nitrobenzene-d5	9.210	82	265727	70.25	ng	0.00
42) 2,4,6-Tribromophenol	16.157	330	96961	70.24	ng	0.00
45) 2-Fluorobiphenyl	13.298	172	598856	72.12	ng	0.00
79) Terphenyl-d14	19.986	244	601416	80.49	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.334	88	34558	32.56	ng	# 67
3) Pyridine	3.764	79	104509	35.56	ng	# 49
4) n-Nitrosodimethylamine	3.675	42	55895	33.08	ng	# 61
6) Aniline	7.352	93	150456	38.55	ng	97
8) 2-Chlorophenol	7.593	128	109649	37.27	ng	96
9) Benzaldehyde	7.163	77	69247	36.94	ng	80
10) Phenol	7.216	94	127289	37.49	ng	87
11) bis(2-Chloroethyl)ether	7.452	93	105217	38.43	ng	91
12) 1,3-Dichlorobenzene	7.916	146	126352	36.40	ng	# 93
13) 1,4-Dichlorobenzene	8.069	146	128536	36.61	ng	98
14) 1,2-Dichlorobenzene	8.387	146	124099	36.85	ng	98
15) Benzyl Alcohol	8.275	79	99251	39.54	ng	# 88
16) 2,2'-oxybis(1-Chloropr...	8.569	45	184334	37.30	ng	71
17) 2-Methylphenol	8.481	107	92250	39.24	ng	# 90
18) Hexachloroethane	9.116	117	48293	36.55	ng	90
19) n-Nitroso-di-n-propyla...	8.851	70	88456	41.23	ng	# 67
20) 3+4-Methylphenols	8.816	107	126805	40.22	ng	91
22) Acetophenone	8.869	105	166913	35.88	ng	# 89
24) Nitrobenzene	9.251	77	128663	35.23	ng	# 88
25) Isophorone	9.775	82	221760	38.26	ng	96
26) 2-Nitrophenol	9.963	139	63441	38.39	ng	# 80
27) 2,4-Dimethylphenol	10.016	122	88589	36.87	ng	91
28) bis(2-Chloroethoxy)met...	10.257	93	151613	37.06	ng	99
29) 2,4-Dichlorophenol	10.493	162	105883	36.93	ng	99
30) 1,2,4-Trichlorobenzene	10.710	180	119968	35.22	ng	98
31) Naphthalene	10.898	128	362418	35.75	ng	99
32) Benzoic acid	10.181	122	87465	40.45	ng	# 86
33) 4-Chloroaniline	11.010	127	154613	36.71	ng	92
34) Hexachlorobutadiene	11.175	225	71754	35.31	ng	98
35) Caprolactam	11.816	113	33024	35.59	ng	# 58
36) 4-Chloro-3-methylphenol	12.128	107	111623	37.44	ng	93
37) 2-Methylnaphthalene	12.504	142	257696	36.88	ng	97
38) 1-Methylnaphthalene	12.722	142	244141	36.82	ng	100
40) 1,2,4,5-Tetrachloroben...	12.869	216	122284	35.59	ng	99
41) Hexachlorocyclopentadiene	12.845	237	61250	38.15	ng	99
43) 2,4,6-Trichlorophenol	13.110	196	86516	37.53	ng	97

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	13.181	196	87529	36.69	ng	97
46) 1,1'-Biphenyl	13.510	154	318485	35.82	ng	99
47) 2-Chloronaphthalene	13.557	162	258780	35.61	ng	98
48) 2-Nitroaniline	13.757	65	80928	35.52	ng	92
49) Acenaphthylene	14.398	152	386985	35.89	ng	99
50) Dimethylphthalate	14.128	163	298179	34.81	ng	99
51) 2,6-Dinitrotoluene	14.251	165	65784	36.10	ng	89
52) Acenaphthene	14.739	154	237426	35.47	ng	98
53) 3-Nitroaniline	14.580	138	71487	34.80	ng	89
54) 2,4-Dinitrophenol	14.786	184	37855	34.88	ng	95
55) Dibenzofuran	15.069	168	357594	34.73	ng	100
56) 4-Nitrophenol	14.880	139	53727	33.22	ng	# 76
57) 2,4-Dinitrotoluene	15.033	165	83980	34.60	ng	93
58) Fluorene	15.716	166	289942	34.37	ng	99
59) 2,3,4,6-Tetrachlorophenol	15.292	232	75363	35.74	ng	94
60) Diethylphthalate	15.480	149	292199	33.60	ng	98
61) 4-Chlorophenyl-phenyle...	15.704	204	143780	34.61	ng	95
62) 4-Nitroaniline	15.739	138	71663	32.05	ng	# 83
63) Azobenzene	15.998	77	280101	34.31	ng	91
65) 4,6-Dinitro-2-methylph...	15.798	198	42857	36.78	ng	81
66) n-Nitrosodiphenylamine	15.922	169	243350	38.75	ng	98
67) 4-Bromophenyl-phenylether	16.592	248	86053	38.74	ng	95
68) Hexachlorobenzene	16.710	284	96978	37.57	ng	96
69) Atrazine	16.857	200	67279	36.14	ng	96
70) Pentachlorophenol	17.051	266	53656	38.14	ng	97
71) Phenanthrene	17.445	178	404967	35.11	ng	99
72) Anthracene	17.533	178	396787	35.62	ng	99
73) Carbazole	17.798	167	350246	33.15	ng	99
74) Di-n-butylphthalate	18.345	149	439172	34.55	ng	99
75) Fluoranthene	19.433	202	399370	30.83	ng	96
77) Benzidine	19.615	184	146823	48.39	ng	100
78) Pyrene	19.792	202	400529	41.40	ng	99
80) Butylbenzylphthalate	20.668	149	168668	39.54	ng	97
81) Benzo(a)anthracene	21.515	228	338390	36.32	ng	100
82) 3,3'-Dichlorobenzidine	21.445	252	113588	37.94	ng	# 98
83) Chrysene	21.562	228	322250	35.83	ng	100
84) Bis(2-ethylhexyl)phtha...	21.427	149	230338	38.19	ng	97
85) Di-n-octyl phthalate	22.315	149	369435	36.23	ng	# 90
87) Indeno(1,2,3-cd)pyrene	26.150	276	347675	37.06	ng	# 90
88) Benzo(b)fluoranthene	23.127	252	327349	38.29	ng	# 97
89) Benzo(k)fluoranthene	23.174	252	299757	35.91	ng	# 97
90) Benzo(a)pyrene	23.715	252	292859	36.97	ng	# 96
91) Dibenzo(a,h)anthracene	26.162	278	288308	37.14	ng	# 94
92) Benzo(g,h,i)perylene	26.868	276	282702	37.03	ng	# 91

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

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