

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP061621\
 Data File : BP005955.D
 Acq On : 16 Jun 2021 18:49
 Operator : CG/JU
 Sample : M2672-21MS
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 17-B6(8-12)MS

Manual Integrations
 APPROVED

mohammad
 6/17/2021 12:16:31 PM

Quant Time: Jun 17 01:58:00 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP060821.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jun 16 11:35:30 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.928	152	71502	20.000	ng	0.00
21) Naphthalene-d8	10.734	136	278268	20.000	ng	0.00
39) Acenaphthene-d10	14.557	164	173210	20.000	ng	0.00
64) Phenanthrene-d10	17.304	188	355970	20.000	ng	0.00
76) Chrysene-d12	21.375	240	426512	20.000	ng	0.00
86) Perylene-d12	23.786	264	497531	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.505	112	593208	133.126	ng	0.00
7) Phenol-d6	7.105	99	639473	110.720	ng	0.00
23) Nitrobenzene-d5	9.099	82	486381	85.694	ng	0.00
42) 2,4,6-Tribromophenol	16.051	330	405302	136.354	ng	0.00
45) 2-Fluorobiphenyl	13.187	172	1120173	84.922	ng	0.00
79) Terphenyl-d14	19.851	244	1982140	87.021	ng	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.376	88	79869	39.674	ng	# 47
3) Pyridine	3.793	79	150160	31.736	ng	99
4) n-Nitrosodimethylamine	3.693	42	87783	39.913	ng	87
6) Aniline	7.258	93	215758	33.461	ng	99
8) 2-Chlorophenol	7.499	128	235926	46.178	ng	98
9) Benzaldehyde	7.069	77	126131	51.228	ng	97
10) Phenol	7.134	94	231953	40.202	ng	98
11) bis(2-Chloroethyl)ether	7.358	93	208855	47.770	ng	98
12) 1,3-Dichlorobenzene	7.816	146	251156	43.963	ng	98
13) 1,4-Dichlorobenzene	7.963	146	259912	44.612	ng	99
14) 1,2-Dichlorobenzene	8.281	146	248424	44.609	ng	98
15) Benzyl Alcohol	8.175	79	204903	47.102	ng	97
16) 2,2'-oxybis(1-Chloropr...	8.458	45	179764	44.479	ng	97
17) 2-Methylphenol	8.381	107	178764	45.474	ng	98
18) Hexachloroethane	9.011	117	93441	44.340	ng	95
19) n-Nitroso-di-n-propyla...	8.746	70	161462	45.443	ng	95
20) 3+4-Methylphenols	8.711	107	235164	44.286	ng	97
22) Acetophenone	8.758	105	338425	46.206	ng	# 98
24) Nitrobenzene	9.140	77	247866	42.055	ng	99
25) Isophorone	9.669	82	425536	40.598	ng	100
26) 2-Nitrophenol	9.852	139	122501	43.562	ng	96
27) 2,4-Dimethylphenol	9.910	122	205666	49.091	ng	98
28) bis(2-Chloroethoxy)met...	10.146	93	267055	48.818	ng	100
29) 2,4-Dichlorophenol	10.393	162	217130	43.127	ng	100
30) 1,2,4-Trichlorobenzene	10.599	180	239779	42.385	ng	97
31) Naphthalene	10.787	128	696617	43.267	ng	100
32) Benzoic acid	10.075	122	112733	37.506	ng	94
33) 4-Chloroaniline	10.905	127	79947	12.292	ng	99
34) Hexachlorobutadiene	11.069	225	158532	41.431	ng	99
35) Caprolactam	11.699	113	53002m	36.548	ng	
36) 4-Chloro-3-methylphenol	12.028	107	228381	44.674	ng	95
37) 2-Methylnaphthalene	12.393	142	503463	43.920	ng	99
38) 1-Methylnaphthalene	12.610	142	465297	43.092	ng	99
40) 1,2,4,5-Tetrachloroben...	12.757	216	269993	44.284	ng	100
41) Hexachlorocyclopentadiene	12.734	237	254566	82.888	ng	98
43) 2,4,6-Trichlorophenol	12.999	196	178211	42.615	ng	96

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	13.075	196	195104	43.321	ng	96
46) 1,1'-Biphenyl	13.399	154	612196	43.836	ng	100
47) 2-Chloronaphthalene	13.440	162	488745	42.857	ng	98
48) 2-Nitroaniline	13.646	65	133148	44.406	ng	94
49) Acenaphthylene	14.281	152	762079	43.558	ng	99
50) Dimethylphthalate	14.022	163	642658	45.126	ng	100
51) 2,6-Dinitrotoluene	14.140	165	140442	43.387	ng	98
52) Acenaphthene	14.622	154	452326	42.572	ng	99
53) 3-Nitroaniline	14.469	138	85227	27.166	ng	99
54) 2,4-Dinitrophenol	14.681	184	139799	73.065	ng	98
55) Dibenzofuran	14.957	168	728972	41.720	ng	98
56) 4-Nitrophenol	14.787	139	201233	79.134	ng	98
57) 2,4-Dinitrotoluene	14.928	165	194206	44.858	ng	# 99
58) Fluorene	15.604	166	588574	43.233	ng	99
59) 2,3,4,6-Tetrachlorophenol	15.187	232	170359m	42.920	ng	
60) Diethylphthalate	15.381	149	648764	45.405	ng	99
61) 4-Chlorophenyl-phenyle...	15.598	204	303130	42.939	ng	98
62) 4-Nitroaniline	15.634	138	137087	42.453	ng	95
63) Azobenzene	15.893	77	538923	41.690	ng	95
65) 4,6-Dinitro-2-methylph...	15.693	198	98846	37.786	ng	97
66) n-Nitrosodiphenylamine	15.816	169	521738	44.198	ng	99
67) 4-Bromophenyl-phenylether	16.492	248	202805	44.388	ng	98
68) Hexachlorobenzene	16.610	284	242890	44.351	ng	98
69) Atrazine	16.769	200	200943	54.714	ng	99
70) Pentachlorophenol	16.957	266	300998	98.846	ng	98
71) Phenanthrene	17.345	178	953168	44.587	ng	99
72) Anthracene	17.439	178	961608	45.705	ng	100
73) Carbazole	17.710	167	881549	43.641	ng	100
74) Di-n-butylphthalate	18.251	149	1139444	48.687	ng	100
75) Fluoranthene	19.304	202	1172597	45.167	ng	98
77) Benzidine	19.486	184	331363	37.337	ng	100
78) Pyrene	19.657	202	1220933	41.005	ng	99
80) Butylbenzylphthalate	20.522	149	561059	45.801	ng	97
81) Benzo(a)anthracene	21.363	228	1305846	42.414	ng	99
82) 3,3'-Dichlorobenzidine	21.292	252	344674	32.381	ng	99
83) Chrysene	21.416	228	1299721	43.585	ng	100
84) Bis(2-ethylhexyl)phtha...	21.280	149	861562	47.956	ng	99
85) Di-n-octyl phthalate	22.198	149	1559426	48.521	ng	99
87) Indeno(1,2,3-cd)pyrene	26.333	276	1875207	44.074	ng	99
88) Benzo(b)fluoranthene	23.051	252	1534758	45.039	ng	100
89) Benzo(k)fluoranthene	23.098	252	1453051	43.974	ng	99
90) Benzo(a)pyrene	23.680	252	1468356	45.833	ng	98
91) Dibenzo(a,h)anthracene	26.339	278	1625554	44.390	ng	99
92) Benzo(g,h,i)perylene	27.110	276	1595062	44.093	ng	99

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

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