

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF061421\
 Data File : BF124309.D
 Acq On : 14 Jun 2021 16:39
 Operator : JU/CG
 Sample : M2626-07MSD
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 18-B12(218-222)MSD

Manual Integrations
 APPROVED

mohammad
 6/15/2021 2:39:13 PM

Quant Time: Jun 14 18:19:12 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF060321.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jun 14 12:10:02 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.828	152	32282	20.000	ng	0.00
21) Naphthalene-d8	8.116	136	115250	20.000	ng	# 0.00
39) Acenaphthene-d10	9.874	164	67462	20.000	ng	0.00
64) Phenanthrene-d10	11.363	188	130918	20.000	ng	0.00
76) Chrysene-d12	14.015	240	80956	20.000	ng	0.00
86) Perylene-d12	15.492	264	70898	20.000	ng	# 0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.481	112	285213	132.998	ng	0.02
7) Phenol-d6	6.486	99	316528	109.809	ng	0.00
23) Nitrobenzene-d5	7.398	82	299530	103.078	ng	0.00
42) 2,4,6-Tribromophenol	10.674	330	141588	148.140	ng	0.00
45) 2-Fluorobiphenyl	9.192	172	558398	104.158	ng	0.00
79) Terphenyl-d14	12.957	244	617320	104.693	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	2.728	88	37132	39.566	ng	# 77
3) Pyridine	3.451	79	67442	28.059	ng	# 81
4) n-Nitrosodimethylamine	3.381	42	62775	43.955	ng	# 81
6) Aniline	6.498	93	52511	15.949	ng	# 1
8) 2-Chlorophenol	6.628	128	112922	47.322	ng	93
9) Benzaldehyde	6.386	77	58981	46.033	ng	92
10) Phenol	6.498	94	121583	39.029	ng	79
11) bis(2-Chloroethyl)ether	6.569	93	119449	52.598	ng	89
12) 1,3-Dichlorobenzene	6.775	146	93689m	33.586	ng	
13) 1,4-Dichlorobenzene	6.851	146	100047	35.755	ng	93
14) 1,2-Dichlorobenzene	6.998	146	96281	35.969	ng	95
15) Benzyl Alcohol	6.981	79	119263	45.896	ng	93
16) 2,2'-oxybis(1-Chloropr...	7.104	45	145081	40.952	ng	# 1
17) 2-Methylphenol	7.098	107	92751	47.815	ng	# 84
18) Hexachloroethane	7.339	117	34453	31.267	ng	# 88
19) n-Nitroso-di-n-propyla...	7.251	70	95961	47.159	ng	# 85
20) 3+4-Methylphenols	7.251	107	117792	45.829	ng	93
22) Acetophenone	7.245	105	178389	54.296	ng	# 95
24) Nitrobenzene	7.422	77	139018	47.689	ng	97
25) Isophorone	7.657	82	238500	45.547	ng	97
26) 2-Nitrophenol	7.733	139	65530	50.323	ng	96
27) 2,4-Dimethylphenol	7.780	122	106414	57.952	ng	91
28) bis(2-Chloroethoxy)met...	7.863	93	139143	53.599	ng	99
29) 2,4-Dichlorophenol	7.986	162	105424	49.618	ng	92
30) 1,2,4-Trichlorobenzene	8.057	180	99392	41.741	ng	97
31) Naphthalene	8.139	128	295452	42.878	ng	99
32) Benzoic acid	7.922	122	49152	43.623	ng	# 80
33) 4-Chloroaniline	8.192	127	10360	4.042	ng	# 90
34) Hexachlorobutadiene	8.251	225	58374	34.707	ng	96
35) Caprolactam	8.580	113	22478m	38.509	ng	
36) 4-Chloro-3-methylphenol	8.686	107	113452	48.805	ng	88
37) 2-Methylnaphthalene	8.827	142	232920	48.203	ng	98
38) 1-Methylnaphthalene	8.927	142	215579	47.742	ng	96
40) 1,2,4,5-Tetrachloroben...	8.998	216	122063	51.068	ng	99
41) Hexachlorocyclopentadiene	8.980	237	78932	62.725	ng	92
43) 2,4,6-Trichlorophenol	9.116	196	77040	49.420	ng	92

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF061421\
 Data File : BF124309.D
 Acq On : 14 Jun 2021 16:39
 Operator : JU/CG
 Sample : M2626-07MSD
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 18-B12(218-222)MSD

Manual Integrations
 APPROVED

mohammad
 6/15/2021 2:39:13 PM

Quant Time: Jun 14 18:19:12 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF060321.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jun 14 12:10:02 2021
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	9.163	196	85121	50.865	ng	95
46) 1,1'-Biphenyl	9.298	154	294811	51.768	ng	98
47) 2-Chloronaphthalene	9.322	162	230350	49.694	ng	98
48) 2-Nitroaniline	9.422	65	91118	54.470	ng	93
49) Acenaphthylene	9.739	152	354865	52.667	ng	98
50) Dimethylphthalate	9.604	163	296938	53.201	ng	99
51) 2,6-Dinitrotoluene	9.663	165	65349	51.064	ng	92
52) Acenaphthene	9.910	154	221588	50.352	ng	95
53) 3-Nitroaniline	9.833	138	32154	26.845	ng	99
54) 2,4-Dinitrophenol	9.951	184	72785	109.869	ng #	88
55) Dibenzofuran	10.080	168	348954	50.663	ng	99
56) 4-Nitrophenol	10.021	139	76198	89.018	ng	89
57) 2,4-Dinitrotoluene	10.074	165	93902	55.636	ng #	89
58) Fluorene	10.427	166	281576	53.258	ng	93
59) 2,3,4,6-Tetrachlorophenol	10.204	232	71861	52.990	ng	93
60) Diethylphthalate	10.298	149	300940	53.426	ng	96
61) 4-Chlorophenyl-phenyle...	10.416	204	139209	51.496	ng	97
62) 4-Nitroaniline	10.457	138	64651	53.683	ng #	78
63) Azobenzene	10.574	77	287894	49.031	ng	96
65) 4,6-Dinitro-2-methylph...	10.486	198	47633	45.274	ng #	74
66) n-Nitrosodiphenylamine	10.539	169	239188	48.085	ng	96
67) 4-Bromophenyl-phenylether	10.910	248	87854	48.891	ng	98
68) Hexachlorobenzene	10.980	284	101370	49.806	ng #	93
69) Atrazine	11.068	200	84653	60.118	ng	95
70) Pentachlorophenol	11.180	266	75554	71.002	ng	94
71) Phenanthrene	11.392	178	412213	50.404	ng	99
72) Anthracene	11.445	178	426338	51.763	ng	98
73) Carbazole	11.604	167	355075	49.471	ng	99
74) Di-n-butylphthalate	11.927	149	461927	50.078	ng	99
75) Fluoranthene	12.586	202	416151	48.477	ng	96
77) Benzidine	12.704	184	37053	22.038	ng	94
78) Pyrene	12.815	202	425632	54.752	ng	98
80) Butylbenzylphthalate	13.427	149	162241	51.447	ng #	93
81) Benzo(a)anthracene	14.004	228	288861	48.980	ng	96
82) 3,3'-Dichlorobenzidine	13.962	252	51483	29.756	ng #	99
83) Chrysene	14.039	228	280237	49.250	ng	97
84) Bis(2-ethylhexyl)phtha...	13.986	149	217903	58.088	ng	98
85) Di-n-octyl phthalate	14.598	149	289319	57.790	ng #	91
87) Indeno(1,2,3-cd)pyrene	16.986	276	305332	51.406	ng	98
88) Benzo(b)fluoranthene	15.062	252	257263	46.712	ng	96
89) Benzo(k)fluoranthene	15.092	252	256875	49.269	ng	95
90) Benzo(a)pyrene	15.433	252	254852	53.276	ng	99
91) Dibenzo(a,h)anthracene	17.003	278	259315	51.192	ng	99
92) Benzo(g,h,i)perylene	17.439	276	268183	53.491	ng	97

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

