

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF061621\
 Data File : BF124339.D
 Acq On : 16 Jun 2021 11:04
 Operator : JU/CG
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDCCC040

Manual Integrations
 APPROVED

mohammad
 6/17/2021 12:15:48 PM

Quant Time: Jun 16 12:19:33 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF060321.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jun 15 20:09:26 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.787	152	38734	20.000	ng	0.00
21) Naphthalene-d8	8.069	136	148779	20.000	ng	0.00
39) Acenaphthene-d10	9.828	164	87958	20.000	ng	0.00
64) Phenanthrene-d10	11.316	188	153113	20.000	ng	0.00
76) Chrysene-d12	13.957	240	83235	20.000	ng	# 0.00
86) Perylene-d12	15.415	264	83594	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.410	112	215095	83.594	ng	0.00
7) Phenol-d6	6.440	99	284535	82.268	ng	0.00
23) Nitrobenzene-d5	7.357	82	296220	78.966	ng	0.00
42) 2,4,6-Tribromophenol	10.622	330	85972	68.990	ng	0.00
45) 2-Fluorobiphenyl	9.151	172	552876	79.097	ng	0.00
79) Terphenyl-d14	12.898	244	466043	76.874	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	2.481	88	44308	39.348	ng	93
3) Pyridine	3.216	79	115013	39.881	ng	92
4) n-Nitrosodimethylamine	3.181	42	66943	39.065	ng	98
6) Aniline	6.457	93	145041	36.715	ng	# 45
8) 2-Chlorophenol	6.581	128	121889	42.572	ng	93
9) Benzaldehyde	6.340	77	52376	34.069	ng	96
10) Phenol	6.451	94	153525	41.074	ng	# 61
11) bis(2-Chloroethyl)ether	6.528	93	116543	42.770	ng	90
12) 1,3-Dichlorobenzene	6.728	146	135659m	40.530	ng	
13) 1,4-Dichlorobenzene	6.804	146	135008	40.213	ng	96
14) 1,2-Dichlorobenzene	6.957	146	129991	40.474	ng	98
15) Benzyl Alcohol	6.939	79	115398	37.011	ng	# 89
16) 2,2'-oxybis(1-Chloropr...	7.063	45	156627	36.846	ng	# 29
17) 2-Methylphenol	7.057	107	97935	42.078	ng	# 84
18) Hexachloroethane	7.298	117	54568	41.272	ng	# 82
19) n-Nitroso-di-n-propyla...	7.210	70	100125	41.009	ng	# 85
20) 3+4-Methylphenols	7.210	107	131749	42.721	ng	# 89
22) Acetophenone	7.204	105	169076	39.864	ng	# 96
24) Nitrobenzene	7.375	77	148265	39.399	ng	93
25) Isophorone	7.616	82	257742	38.129	ng	96
26) 2-Nitrophenol	7.692	139	74374	44.243	ng	# 86
27) 2,4-Dimethylphenol	7.734	122	95189	40.157	ng	94
28) bis(2-Chloroethoxy)met...	7.822	93	134315	40.079	ng	97
29) 2,4-Dichlorophenol	7.939	162	112920	41.169	ng	91
30) 1,2,4-Trichlorobenzene	8.010	180	121139	39.409	ng	92
31) Naphthalene	8.092	128	349218	39.260	ng	98
32) Benzoic acid	7.881	122	53365	36.689	ng	94
33) 4-Chloroaniline	8.151	127	114726	34.676	ng	96
34) Hexachlorobutadiene	8.210	225	84110	38.739	ng	97
35) Caprolactam	8.528	113	31292m	41.527	ng	
36) 4-Chloro-3-methylphenol	8.639	107	115631	38.532	ng	93
37) 2-Methylnaphthalene	8.786	142	245672	39.384	ng	98
38) 1-Methylnaphthalene	8.886	142	230178	39.487	ng	93
40) 1,2,4,5-Tetrachloroben...	8.951	216	123481	39.623	ng	# 96
41) Hexachlorocyclopentadiene	8.939	237	58419	38.938	ng	97
43) 2,4,6-Trichlorophenol	9.069	196	77293	38.028	ng	95

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF061621\
 Data File : BF124339.D
 Acq On : 16 Jun 2021 11:04
 Operator : JU/CG
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDCCC040

Manual Integrations
 APPROVED

mohammad
 6/17/2021 12:15:48 PM

Quant Time: Jun 16 12:19:33 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF060321.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jun 15 20:09:26 2021
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	9.122	196	85362	39.123	ng	# 91
46) 1,1'-Biphenyl	9.251	154	297377	40.051	ng	98
47) 2-Chloronaphthalene	9.275	162	236257	39.092	ng	96
48) 2-Nitroaniline	9.380	65	83769	38.408	ng	92
49) Acenaphthylene	9.692	152	350743	39.925	ng	98
50) Dimethylphthalate	9.557	163	274667	37.744	ng	99
51) 2,6-Dinitrotoluene	9.622	165	65066	38.995	ng	# 84
52) Acenaphthene	9.863	154	223577	38.966	ng	94
53) 3-Nitroaniline	9.792	138	60253	38.583	ng	# 93
54) 2,4-Dinitrophenol	9.910	184	28321m	35.252	ng	
55) Dibenzofuran	10.033	168	351334	39.123	ng	98
56) 4-Nitrophenol	9.975	139	44703	40.055	ng	82
57) 2,4-Dinitrotoluene	10.028	165	87341	39.690	ng	# 89
58) Fluorene	10.380	166	269265	39.062	ng	94
59) 2,3,4,6-Tetrachlorophenol	10.157	232	64174	36.295	ng	92
60) Diethylphthalate	10.251	149	278588	37.933	ng	99
61) 4-Chlorophenyl-phenyle...	10.369	204	131515	37.313	ng	96
62) 4-Nitroaniline	10.410	138	50281	32.022	ng	# 83
63) Azobenzene	10.527	77	283465	37.027	ng	97
65) 4,6-Dinitro-2-methylph...	10.439	198	41470	33.703	ng	# 69
66) n-Nitrosodiphenylamine	10.492	169	230508	39.622	ng	98
67) 4-Bromophenyl-phenylether	10.857	248	82589	39.298	ng	95
68) Hexachlorobenzene	10.927	284	90621	38.071	ng	94
69) Atrazine	11.022	200	56534	34.329	ng	93
70) Pentachlorophenol	11.127	266	37432m	32.766	ng	
71) Phenanthrene	11.339	178	373068	39.005	ng	98
72) Anthracene	11.392	178	377397	39.179	ng	98
73) Carbazole	11.551	167	315863	37.628	ng	96
74) Di-n-butylphthalate	11.874	149	368540	34.162	ng	98
75) Fluoranthene	12.527	202	342229	34.087	ng	97
77) Benzidine	12.657	184	33173	19.715	ng	# 85
78) Pyrene	12.757	202	331858	41.520	ng	100
80) Butylbenzylphthalate	13.374	149	115315	35.566	ng	98
81) Benzo(a)anthracene	13.951	228	244307	40.291	ng	98
82) 3,3'-Dichlorobenzidine	13.910	252	69435	39.033	ng	96
83) Chrysene	13.986	228	236819	40.480	ng	96
84) Bis(2-ethylhexyl)phtha...	13.933	149	140924	36.538	ng	97
85) Di-n-octyl phthalate	14.545	149	243569	47.320	ng	# 91
87) Indeno(1,2,3-cd)pyrene	16.874	276	280636	41.028	ng	96
88) Benzo(b)fluoranthene	14.998	252	252543	38.891	ng	97
89) Benzo(k)fluoranthene	15.021	252	228345	37.145	ng	99
90) Benzo(a)pyrene	15.357	252	226241	40.112	ng	97
91) Dibenzo(a,h)anthracene	16.880	278	234538	40.241	ng	98
92) Benzo(g,h,i)perylene	17.309	276	233475	40.492	ng	98

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

