

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF111220\
 Data File : BF122312.D
 Acq On : 12 Nov 2020 13:07
 Operator : JU/CG
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDCCC040

Manual Integrations
 APPROVED

mohammad
 11/20/2020 11:27:50 AM

Quant Time: Nov 12 13:40:18 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF110920.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Nov 10 12:25:42 2020
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.869	152	292305	20.00	ng	0.00
21) Naphthalene-d8	8.151	136	1066314	20.00	ng	0.00
39) Acenaphthene-d10	9.910	164	556335	20.00	ng	0.00
64) Phenanthrene-d10	11.392	188	992306	20.00	ng	0.00
76) Chrysene-d12	14.033	240	846901	20.00	ng	0.00
86) Perylene-d12	15.498	264	744600	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.475	112	1500476	78.82	ng	0.00
7) Phenol-d6	6.493	99	1937383	77.80	ng	0.00
23) Nitrobenzene-d5	7.434	82	1543285	88.60	ng	0.00
42) 2,4,6-Tribromophenol	10.698	330	495272	82.95	ng	0.00
45) 2-Fluorobiphenyl	9.234	172	2670973	75.01	ng	0.00
79) Terphenyl-d14	12.980	244	3048479	76.26	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	2.652	88	339507	38.89	ng	93
3) Pyridine	3.387	79	934508	38.92	ng	93
4) n-Nitrosodimethylamine	3.346	42	439933	38.87	ng	99
6) Aniline	6.528	93	1266049	38.75	ng	98
8) 2-Chlorophenol	6.651	128	786575	39.84	ng	92
9) Benzaldehyde	6.416	77	508914	38.00	ng	96
10) Phenol	6.504	94	967764m	39.46	ng	
11) bis(2-Chloroethyl)ether	6.604	93	759548	38.97	ng	94
12) 1,3-Dichlorobenzene	6.810	146	883141	39.40	ng	97
13) 1,4-Dichlorobenzene	6.887	146	879933	39.08	ng	98
14) 1,2-Dichlorobenzene	7.040	146	833799	39.68	ng	99
15) Benzyl Alcohol	7.004	79	693673	39.18	ng	99
16) 2,2'-oxybis(1-Chloropr...	7.145	45	1091284	38.41	ng	98
17) 2-Methylphenol	7.110	107	645923	39.20	ng	99
18) Hexachloroethane	7.387	117	314935	40.17	ng	95
19) n-Nitroso-di-n-propyla...	7.287	70	569005	38.23	ng	95
20) 3+4-Methylphenols	7.269	107	798425	39.38	ng	# 74
22) Acetophenone	7.281	105	1065254	38.35	ng	# 84
24) Nitrobenzene	7.451	77	837709	42.04	ng	96
25) Isophorone	7.692	82	1478391	38.07	ng	97
26) 2-Nitrophenol	7.769	139	335411	40.08	ng	96
27) 2,4-Dimethylphenol	7.798	122	740702	40.44	ng	99
28) bis(2-Chloroethoxy)met...	7.904	93	921223	38.79	ng	97
29) 2,4-Dichlorophenol	8.004	162	659463	40.67	ng	97
30) 1,2,4-Trichlorobenzene	8.092	180	709627	39.65	ng	96
31) Naphthalene	8.175	128	2196981	38.75	ng	98
32) Benzoic acid	7.910	122	371940m	38.56	ng	
33) 4-Chloroaniline	8.222	127	973180	39.42	ng	95
34) Hexachlorobutadiene	8.298	225	408931	39.75	ng	99
35) Caprolactam	8.586	113	204423	39.77	ng	91
36) 4-Chloro-3-methylphenol	8.698	107	671489	39.70	ng	91
37) 2-Methylnaphthalene	8.869	142	1451063	38.74	ng	99
38) 1-Methylnaphthalene	8.969	142	1360295	38.80	ng	99
40) 1,2,4,5-Tetrachloroben...	9.034	216	678583	39.58	ng	99
41) Hexachlorocyclopentadiene	9.022	237	419369	41.84	ng	97
43) 2,4,6-Trichlorophenol	9.139	196	453856	40.76	ng	99

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF111220\
 Data File : BF122312.D
 Acq On : 12 Nov 2020 13:07
 Operator : JU/CG
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDCCC040

Manual Integrations
 APPROVED

mohammad
 11/20/2020 11:27:50 AM

Quant Time: Nov 12 13:40:18 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF110920.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Nov 10 12:25:42 2020
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	9.175	196	473479	40.57	ng	99
46) 1,1'-Biphenyl	9.334	154	1714031	38.71	ng	98
47) 2-Chloronaphthalene	9.357	162	1373277	39.07	ng	97
48) 2-Nitroaniline	9.445	65	411213	40.21	ng	96
49) Acenaphthylene	9.769	152	2116133	39.17	ng	98
50) Dimethylphthalate	9.633	163	1536859	38.86	ng	100
51) 2,6-Dinitrotoluene	9.686	165	350315	40.46	ng	97
52) Acenaphthene	9.945	154	1310128	39.18	ng	99
53) 3-Nitroaniline	9.857	138	391104	39.68	ng	96
54) 2,4-Dinitrophenol	9.957	184	116477	45.74	ng	# 83
55) Dibenzofuran	10.116	168	1904495	38.88	ng	99
56) 4-Nitrophenol	10.004	139	314573	39.41	ng	96
57) 2,4-Dinitrotoluene	10.092	165	452704	41.53	ng	# 85
58) Fluorene	10.457	166	1426074	38.02	ng	100
59) 2,3,4,6-Tetrachlorophenol	10.228	232	402711	41.49	ng	99
60) Diethylphthalate	10.333	149	1510327	38.93	ng	99
61) 4-Chlorophenyl-phenyle...	10.451	204	687785	38.84	ng	98
62) 4-Nitroaniline	10.469	138	380005	39.40	ng	92
63) Azobenzene	10.610	77	1551658	38.13	ng	97
65) 4,6-Dinitro-2-methylph...	10.498	198	183087	44.75	ng	95
66) n-Nitrosodiphenylamine	10.569	169	1322136	39.11	ng	95
67) 4-Bromophenyl-phenylether	10.939	248	464659	39.85	ng	99
68) Hexachlorobenzene	11.004	284	500461	39.72	ng	98
69) Atrazine	11.092	200	333081	39.56	ng	97
70) Pentachlorophenol	11.192	266	300174	41.36	ng	96
71) Phenanthrene	11.416	178	2177320	38.67	ng	97
72) Anthracene	11.469	178	2227167	38.38	ng	98
73) Carbazole	11.622	167	2005946	38.00	ng	99
74) Di-n-butylphthalate	11.957	149	2183663	38.80	ng	99
75) Fluoranthene	12.604	202	2257151	38.40	ng	98
77) Benzidine	12.722	184	883912	37.63	ng	98
78) Pyrene	12.833	202	2337802	39.29	ng	98
80) Butylbenzylphthalate	13.457	149	902454	38.52	ng	93
81) Benzo(a)anthracene	14.021	228	2063008	39.17	ng	99
82) 3,3'-Dichlorobenzidine	13.986	252	776875	39.58	ng	98
83) Chrysene	14.063	228	2070006	39.02	ng	98
84) Bis(2-ethylhexyl)phtha...	14.016	149	1149498	40.07	ng	99
85) Di-n-octyl phthalate	14.633	149	1956632	40.24	ng	99
87) Indeno(1,2,3-cd)pyrene	16.968	276	1676022	37.51	ng	99
88) Benzo(b)fluoranthene	15.074	252	1947250	40.38	ng	99
89) Benzo(k)fluoranthene	15.104	252	1920372	40.84	ng	98
90) Benzo(a)pyrene	15.439	252	1692418	40.26	ng	98
91) Dibenzo(a,h)anthracene	16.986	278	1398611	37.37	ng	99
92) Benzo(g,h,i)perylene	17.409	276	1359594	37.36	ng	98

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

