

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF100322\
 Data File : BF130493.D
 Acq On : 03 Oct 2022 12:21
 Operator : CG\JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDCCC040

Quant Time: Oct 04 03:12:48 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF091422.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Sep 28 02:59:31 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.640	152	84383	20.000	ng	0.00
21) Naphthalene-d8	7.922	136	339586	20.000	ng	0.00
39) Acenaphthene-d10	9.669	164	175244	20.000	ng	-0.01
64) Phenanthrene-d10	11.151	188	345155	20.000	ng	# 0.00
76) Chrysene-d12	13.786	240	162300	20.000	ng	0.00
86) Perylene-d12	15.168	264	141934	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.234	112	392933	80.766	ng	-0.01
7) Phenol-d6	6.287	99	519571	85.353	ng	0.00
23) Nitrobenzene-d5	7.210	82	583817	87.831	ng	0.00
42) 2,4,6-Tribromophenol	10.463	330	134950	80.367	ng	0.00
45) 2-Fluorobiphenyl	8.998	172	951488	79.064	ng	-0.01
79) Terphenyl-d14	12.739	244	817122	83.398	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	2.240	88	146375	65.511	ng	94
3) Pyridine	2.910	79	298531	51.219	ng	97
4) n-Nitrosodimethylamine	2.875	42	154414	48.710	ng	94
6) Aniline	6.304	93	319571	42.607	ng	# 82
8) 2-Chlorophenol	6.422	128	227291	41.013	ng	95
9) Benzaldehyde	6.187	77	166194	40.758	ng	98
10) Phenol	6.298	94	308546	43.252	ng	# 69
11) bis(2-Chloroethyl)ether	6.381	93	219158	42.592	ng	98
12) 1,3-Dichlorobenzene	6.575	146	249063	40.843	ng	96
13) 1,4-Dichlorobenzene	6.657	146	257416	41.395	ng	99
14) 1,2-Dichlorobenzene	6.810	146	244635	42.113	ng	99
15) Benzyl Alcohol	6.792	79	197234	40.866	ng	98
16) 2,2'-oxybis(1-Chloropr...	6.922	45	352493	46.969	ng	96
17) 2-Methylphenol	6.904	107	196048	43.517	ng	96
18) Hexachloroethane	7.145	117	104185	42.503	ng	95
19) n-Nitroso-di-n-propyla...	7.063	70	187523	45.653	ng	98
20) 3+4-Methylphenols	7.063	107	255910	43.728	ng	# 72
22) Acetophenone	7.057	105	350693	42.240	ng	# 99
24) Nitrobenzene	7.228	77	292639	43.358	ng	98
25) Isophorone	7.469	82	472936	44.144	ng	98
26) 2-Nitrophenol	7.545	139	111145	40.171	ng	94
27) 2,4-Dimethylphenol	7.587	122	173577	40.745	ng	99
28) bis(2-Chloroethoxy)met...	7.681	93	268287	44.473	ng	98
29) 2,4-Dichlorophenol	7.787	162	191838	41.093	ng	96
30) 1,2,4-Trichlorobenzene	7.863	180	208416	41.293	ng	98
31) Naphthalene	7.945	128	752389	43.006	ng	99
32) Benzoic acid	7.728	122	97290	29.531	ng	97
33) 4-Chloroaniline	7.998	127	292421	43.947	ng	96
34) Hexachlorobutadiene	8.057	225	126693	39.277	ng	98
35) Caprolactam	8.381	113	63944	47.779	ng	95
36) 4-Chloro-3-methylphenol	8.486	107	238400	46.072	ng	94
37) 2-Methylnaphthalene	8.634	142	540358	48.442	ng	100
38) 1-Methylnaphthalene	8.733	142	457265	42.672	ng	100
40) 1,2,4,5-Tetrachloroben...	8.798	216	185496	38.786	ng	98
41) Hexachlorocyclopentadiene	8.781	237	91637	35.789	ng	97
43) 2,4,6-Trichlorophenol	8.916	196	122949	36.836	ng	96

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF100322\
 Data File : BF130493.D
 Acq On : 03 Oct 2022 12:21
 Operator : CG\JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDCCC040

Quant Time: Oct 04 03:12:48 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF091422.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Sep 28 02:59:31 2022
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	8.957	196	133254	36.997	ng	# 89
46) 1,1'-Biphenyl	9.098	154	533282	40.746	ng	100
47) 2-Chloronaphthalene	9.122	162	431326	40.740	ng	100
48) 2-Nitroaniline	9.228	65	158396	44.856	ng	99
49) Acenaphthylene	9.533	152	665820	41.944	ng	99
50) Dimethylphthalate	9.404	163	498174	41.154	ng	99
51) 2,6-Dinitrotoluene	9.469	165	108557	42.880	ng	99
52) Acenaphthene	9.704	154	402650	38.606	ng	98
53) 3-Nitroaniline	9.639	138	122002	43.248	ng	98
54) 2,4-Dinitrophenol	9.745	184	40814	33.650	ng	93
55) Dibenzofuran	9.875	168	617714	42.580	ng	96
56) 4-Nitrophenol	9.810	139	69223	33.691	ng	97
57) 2,4-Dinitrotoluene	9.869	165	144455	44.647	ng	# 84
58) Fluorene	10.222	166	546859	46.093	ng	100
59) 2,3,4,6-Tetrachlorophenol	9.998	232	105276	37.343	ng	95
60) Diethylphthalate	10.104	149	550346	45.337	ng	100
61) 4-Chlorophenyl-phenyle...	10.216	204	237917	45.713	ng	99
62) 4-Nitroaniline	10.251	138	144176	50.876	ng	93
63) Azobenzene	10.375	77	672170	53.775	ng	99
65) 4,6-Dinitro-2-methylph...	10.275	198	67283	35.018	ng	79
66) n-Nitrosodiphenylamine	10.333	169	482445	41.828	ng	99
67) 4-Bromophenyl-phenylether	10.704	248	155582	40.861	ng	96
68) Hexachlorobenzene	10.763	284	147519	34.178	ng	97
69) Atrazine	10.869	200	130539	38.728	ng	98
70) Pentachlorophenol	10.963	266	78800	34.018	ng	97
71) Phenanthrene	11.180	178	728906	37.614	ng	99
72) Anthracene	11.227	178	763473	40.826	ng	100
73) Carbazole	11.386	167	563780	33.405	ng	98
74) Di-n-butylphthalate	11.722	149	671555	32.997	ng	99
75) Fluoranthene	12.357	202	596888	31.876	ng	98
77) Benzidine	12.492	184	141674	30.070	ng	99
78) Pyrene	12.586	202	604042	42.544	ng	100
80) Butylbenzylphthalate	13.216	149	222144	42.234	ng	96
81) Benzo(a)anthracene	13.774	228	441456	40.887	ng	99
82) 3,3'-Dichlorobenzidine	13.745	252	100213	34.097	ng	98
83) Chrysene	13.810	228	451006	43.993	ng	100
84) Bis(2-ethylhexyl)phtha...	13.774	149	249620	41.432	ng	99
85) Di-n-octyl phthalate	14.386	149	289124	31.376	ng	98
87) Indeno(1,2,3-cd)pyrene	16.492	276	370734	36.834	ng	99
88) Benzo(b)fluoranthene	14.786	252	318971	34.431	ng	99
89) Benzo(k)fluoranthene	14.810	252	323530	35.502	ng	100
90) Benzo(a)pyrene	15.115	252	324287	44.185	ng	99
91) Dibenzo(a,h)anthracene	16.509	278	306248	37.090	ng	98
92) Benzo(g,h,i)perylene	16.892	276	305076	36.678	ng	97

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

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF100322\
 Data File : BF130493.D
 Acq On : 03 Oct 2022 12:21
 Operator : CG\JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDCCC040

Quant Time: Oct 04 03:12:48 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF091422.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Sep 28 02:59:31 2022
 Response via : Initial Calibration

