

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF010825\
 Data File : BF141079.D
 Acq On : 08 Jan 2025 10:49
 Operator : RC/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDCCC040

Quant Time: Jan 08 11:10:43 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF122624.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Dec 27 00:31:16 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.834	152	229689	20.000	ng	0.00
21) Naphthalene-d8	8.116	136	890442	20.000	ng	0.00
39) Acenaphthene-d10	9.863	164	476495	20.000	ng	0.00
64) Phenanthrene-d10	11.351	188	788190	20.000	ng	0.00
76) Chrysene-d12	13.998	240	453606	20.000	ng	0.01
86) Perylene-d12	15.492	264	453696	20.000	ng	0.05
System Monitoring Compounds						
5) 2-Fluorophenol	5.434	112	1130093	74.165	ng	0.00
7) Phenol-d6	6.457	99	1433304	73.018	ng	0.00
23) Nitrobenzene-d5	7.392	82	1283581	92.225	ng	0.00
42) 2,4,6-Tribromophenol	10.651	330	418770	90.445	ng	0.00
45) 2-Fluorobiphenyl	9.186	172	2278209	76.176	ng	0.00
79) Terphenyl-d14	12.939	244	2294557	84.021	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	2.552	88	250929	36.374	ng	97
3) Pyridine	3.299	79	637796	37.831	ng	98
4) n-Nitrosodimethylamine	3.246	42	305271	35.737	ng	99
6) Aniline	6.492	93	667482	38.128	ng	99
8) 2-Chlorophenol	6.616	128	610627	38.009	ng	97
9) Benzaldehyde	6.381	77	377479	39.164	ng	98
10) Phenol	6.469	94	769183	37.896	ng	98
11) bis(2-Chloroethyl)ether	6.569	93	576510	35.177	ng	97
12) 1,3-Dichlorobenzene	6.769	146	675290	38.088	ng	100
13) 1,4-Dichlorobenzene	6.851	146	681437	38.126	ng	99
14) 1,2-Dichlorobenzene	7.004	146	635987	38.165	ng	99
15) Benzyl Alcohol	6.969	79	524966	37.216	ng	99
16) 2,2'-oxybis(1-Chloropr...	7.104	45	857826	34.611	ng	99
17) 2-Methylphenol	7.081	107	494158	37.025	ng	100
18) Hexachloroethane	7.345	117	248407	39.124	ng	99
19) n-Nitroso-di-n-propyla...	7.245	70	424321	35.285	ng	98
20) 3+4-Methylphenols	7.234	107	622639	36.865	ng	# 86
22) Acetophenone	7.239	105	798811	36.569	ng	98
24) Nitrobenzene	7.410	77	674839	43.393	ng	98
25) Isophorone	7.651	82	1127296	37.221	ng	100
26) 2-Nitrophenol	7.728	139	304354	47.782	ng	98
27) 2,4-Dimethylphenol	7.763	122	413750	37.380	ng	100
28) bis(2-Chloroethoxy)met...	7.863	93	710566	37.030	ng	99
29) 2,4-Dichlorophenol	7.963	162	496646	39.562	ng	99
30) 1,2,4-Trichlorobenzene	8.051	180	555284	39.081	ng	100
31) Naphthalene	8.134	128	1795871	38.274	ng	99
32) Benzoic acid	7.869	122	411236	45.604	ng	97
33) 4-Chloroaniline	8.181	127	631644	37.211	ng	99
34) Hexachlorobutadiene	8.251	225	337151	40.250	ng	99
35) Caprolactam	8.551	113	166859	39.099	ng	91
36) 4-Chloro-3-methylphenol	8.657	107	543391	38.977	ng	99
37) 2-Methylnaphthalene	8.822	142	1144198	38.118	ng	100
38) 1-Methylnaphthalene	8.922	142	1124960	38.134	ng	100
40) 1,2,4,5-Tetrachloroben...	8.986	216	519348	39.618	ng	99
41) Hexachlorocyclopentadiene	8.975	237	191303	44.037	ng	98
43) 2,4,6-Trichlorophenol	9.098	196	353659	40.535	ng	99

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF010825\
 Data File : BF141079.D
 Acq On : 08 Jan 2025 10:49
 Operator : RC/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDCCC040

Quant Time: Jan 08 11:10:43 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF122624.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Dec 27 00:31:16 2024
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	9.133	196	380571	42.631	ng	99
46) 1,1'-Biphenyl	9.286	154	1399088	38.347	ng	99
47) 2-Chloronaphthalene	9.310	162	1096589	38.753	ng	100
48) 2-Nitroaniline	9.404	65	331512	45.095	ng	91
49) Acenaphthylene	9.728	152	1563704	38.388	ng	100
50) Dimethylphthalate	9.586	163	1224426	39.066	ng	100
51) 2,6-Dinitrotoluene	9.645	165	289566	45.508	ng	97
52) Acenaphthene	9.898	154	1086234	39.444	ng	100
53) 3-Nitroaniline	9.816	138	300556	43.562	ng	# 92
54) 2,4-Dinitrophenol	9.922	184	130076	59.075	ng	# 1
55) Dibenzofuran	10.069	168	1487133	38.426	ng	99
56) 4-Nitrophenol	9.963	139	241207	44.990	ng	97
57) 2,4-Dinitrotoluene	10.051	165	376001	48.065	ng	94
58) Fluorene	10.416	166	1137869	37.915	ng	100
59) 2,3,4,6-Tetrachlorophenol	10.186	232	318556	43.087	ng	97
60) Diethylphthalate	10.286	149	1198691	38.740	ng	100
61) 4-Chlorophenyl-phenyle...	10.404	204	563083	38.993	ng	99
62) 4-Nitroaniline	10.428	138	291967	42.459	ng	95
63) Azobenzene	10.569	77	1206641	37.153	ng	98
65) 4,6-Dinitro-2-methylph...	10.457	198	189388	57.878	ng	94
66) n-Nitrosodiphenylamine	10.522	169	1010230	40.702	ng	99
67) 4-Bromophenyl-phenylether	10.898	248	370235	43.452	ng	96
68) Hexachlorobenzene	10.963	284	398693	42.436	ng	98
69) Atrazine	11.045	200	239580	34.552	ng	99
70) Pentachlorophenol	11.151	266	270973	46.324	ng	99
71) Phenanthrene	11.375	178	1651051	39.800	ng	100
72) Anthracene	11.427	178	1625433	39.988	ng	99
73) Carbazole	11.580	167	1500292	38.394	ng	99
74) Di-n-butylphthalate	11.916	149	1766742	39.751	ng	99
75) Fluoranthene	12.563	202	1673369	37.193	ng	98
77) Benzidine	12.680	184	422736	44.097	ng	99
78) Pyrene	12.792	202	1645147	41.469	ng	100
80) Butylbenzylphthalate	13.416	149	579093	39.976	ng	98
81) Benzo(a)anthracene	13.992	228	1180479	37.341	ng	99
82) 3,3'-Dichlorobenzidine	13.951	252	382305	40.250	ng	99
83) Chrysene	14.027	228	1061204	37.237	ng	100
84) Bis(2-ethylhexyl)phtha...	13.986	149	691740	36.813	ng	100
85) Di-n-octyl phthalate	14.627	149	1065063	39.162	ng	98
87) Indeno(1,2,3-cd)pyrene	16.957	276	1297695	45.400	ng	97
88) Benzo(b)fluoranthene	15.063	252	1121970	35.458	ng	99
89) Benzo(k)fluoranthene	15.092	252	934620	37.686	ng	99
90) Benzo(a)pyrene	15.427	252	942214	38.363	ng	99
91) Dibenzo(a,h)anthracene	16.980	278	1060367	45.764	ng	98
92) Benzo(g,h,i)perylene	17.404	276	1108046	46.107	ng	98

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

Quant Time: Jan 08 11:10:43 2025
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF122624.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Dec 27 00:31:16 2024
Response via : Initial Calibration

