

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM032023\
 Data File : BM039037.D
 Acq On : 20 Mar 2023 10:24
 Operator : CG/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTDCCC040

Quant Time: Mar 21 04:03:58 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM030823.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Mar 17 05:25:34 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.975	152	358697	20.000	ng	0.00
21) Naphthalene-d8	10.786	136	1449806	20.000	ng	0.00
39) Acenaphthene-d10	14.615	164	805611	20.000	ng	0.00
64) Phenanthrene-d10	17.356	188	1568589	20.000	ng	0.00
76) Chrysene-d12	21.539	240	1372955	20.000	ng	0.00
86) Perylene-d12	23.974	264	1453669	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.522	112	1882126	79.719	ng	0.00
7) Phenol-d6	7.134	99	2480490	80.886	ng	0.00
23) Nitrobenzene-d5	9.145	82	2420704	82.016	ng	0.00
42) 2,4,6-Tribromophenol	16.098	330	711539	76.348	ng	0.00
45) 2-Fluorobiphenyl	13.239	172	4760350	76.923	ng	0.00
79) Terphenyl-d14	19.980	244	6015503	80.107	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.393	88	404833	38.543	ng	97
3) Pyridine	3.799	79	1193636	40.976	ng	97
4) n-Nitrosodimethylamine	3.710	42	496687	41.901	ng	98
6) Aniline	7.298	93	1659985	41.226	ng	99
8) 2-Chlorophenol	7.534	128	1010255	40.193	ng	100
9) Benzaldehyde	7.110	77	555613	38.074	ng	99
10) Phenol	7.157	94	1311396	40.269	ng	100
11) bis(2-Chloroethyl)ether	7.392	93	1080229	40.799	ng	97
12) 1,3-Dichlorobenzene	7.863	146	1043780	39.076	ng	100
13) 1,4-Dichlorobenzene	8.010	146	1056308	38.882	ng	100
14) 1,2-Dichlorobenzene	8.328	146	1019334	38.953	ng	99
15) Benzyl Alcohol	8.216	79	924858	42.077	ng	98
16) 2,2'-oxybis(1-Chloropr...	8.504	45	1453536	43.468	ng	98
17) 2-Methylphenol	8.416	107	921520	41.193	ng	100
18) Hexachloroethane	9.057	117	405831	40.489	ng	99
19) n-Nitroso-di-n-propyla...	8.786	70	865260	44.796	ng	97
20) 3+4-Methylphenols	8.745	107	1244085	41.649	ng	99
22) Acetophenone	8.804	105	1591801	39.133	ng	# 99
24) Nitrobenzene	9.186	77	1252479	40.374	ng	100
25) Isophorone	9.710	82	2284644	42.064	ng	99
26) 2-Nitrophenol	9.898	139	532818	39.465	ng	97
27) 2,4-Dimethylphenol	9.957	122	949972	40.141	ng	98
28) bis(2-Chloroethoxy)met...	10.198	93	1413942	40.516	ng	99
29) 2,4-Dichlorophenol	10.428	162	883696	39.386	ng	99
30) 1,2,4-Trichlorobenzene	10.645	180	911032	37.157	ng	99
31) Naphthalene	10.839	128	3198124	38.685	ng	100
32) Benzoic acid	10.092	122	679038	40.169	ng	100
33) 4-Chloroaniline	10.945	127	1412866	40.102	ng	99
34) Hexachlorobutadiene	11.122	225	514026	36.558	ng	99
35) Caprolactam	11.739	113	284270	41.628	ng	96
36) 4-Chloro-3-methylphenol	12.069	107	986804	41.303	ng	99
37) 2-Methylnaphthalene	12.445	142	2186488	39.460	ng	100
38) 1-Methylnaphthalene	12.663	142	2045525	39.393	ng	100
40) 1,2,4,5-Tetrachloroben...	12.810	216	965999	36.884	ng	99
41) Hexachlorocyclopentadiene	12.792	237	571780	39.773	ng	99
43) 2,4,6-Trichlorophenol	13.051	196	666542	39.302	ng	100

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM032023\
 Data File : BM039037.D
 Acq On : 20 Mar 2023 10:24
 Operator : CG/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTDCCC040

Quant Time: Mar 21 04:03:58 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM030823.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Mar 17 05:25:34 2023
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	13.116	196	775111	39.046	ng	98
46) 1,1'-Biphenyl	13.451	154	2635390	38.614	ng	98
47) 2-Chloronaphthalene	13.492	162	1975056	38.377	ng	100
48) 2-Nitroaniline	13.692	65	679055	41.524	ng	96
49) Acenaphthylene	14.333	152	3260743	39.777	ng	99
50) Dimethylphthalate	14.074	163	2463690	39.743	ng	100
51) 2,6-Dinitrotoluene	14.192	165	532842	39.149	ng	97
52) Acenaphthene	14.680	154	1957108	38.847	ng	100
53) 3-Nitroaniline	14.521	138	628128	39.484	ng	96
54) 2,4-Dinitrophenol	14.721	184	297956	40.549	ng	99
55) Dibenzofuran	15.010	168	3056181	38.766	ng	99
56) 4-Nitrophenol	14.815	139	501888	40.765	ng	96
57) 2,4-Dinitrotoluene	14.974	165	701130	39.255	ng	96
58) Fluorene	15.657	166	2431816	39.379	ng	99
59) 2,3,4,6-Tetrachlorophenol	15.233	232	598036	39.063	ng	98
60) Diethylphthalate	15.433	149	2494434	41.355	ng	99
61) 4-Chlorophenyl-phenyle...	15.651	204	1172383	38.643	ng	97
62) 4-Nitroaniline	15.680	138	643477	39.758	ng	100
63) Azobenzene	15.945	77	2735522	42.861	ng	99
65) 4,6-Dinitro-2-methylph...	15.733	198	393342	39.201	ng	95
66) n-Nitrosodiphenylamine	15.868	169	2070623	39.157	ng	99
67) 4-Bromophenyl-phenylether	16.545	248	657064	38.782	ng	94
68) Hexachlorobenzene	16.662	284	709745	37.473	ng	98
69) Atrazine	16.821	200	592464	41.018	ng	99
70) Pentachlorophenol	17.004	266	537925	38.615	ng	99
71) Phenanthrene	17.398	178	3555059	38.616	ng	100
72) Anthracene	17.492	178	3616829	39.395	ng	100
73) Carbazole	17.762	167	3358382	40.081	ng	99
74) Di-n-butylphthalate	18.327	149	4107219	40.482	ng	100
75) Fluoranthene	19.409	202	3865888	39.771	ng	98
77) Benzidine	19.598	184	1323306	42.865	ng	98
78) Pyrene	19.774	202	4084205	39.606	ng	100
80) Butylbenzylphthalate	20.674	149	1811082	41.767	ng	100
81) Benzo(a)anthracene	21.521	228	3936283	39.811	ng	100
82) 3,3'-Dichlorobenzidine	21.456	252	1275932	41.989	ng	100
83) Chrysene	21.574	228	3756568	39.065	ng	100
84) Bis(2-ethylhexyl)phtha...	21.450	149	2732690	43.358	ng	99
85) Di-n-octyl phthalate	22.386	149	4571975	43.816	ng	99
87) Indeno(1,2,3-cd)pyrene	26.515	276	4372114	39.526	ng	98
88) Benzo(b)fluoranthene	23.233	252	3733788	40.204	ng	99
89) Benzo(k)fluoranthene	23.280	252	3751171	38.812	ng	99
90) Benzo(a)pyrene	23.868	252	3604812	40.331	ng	100
91) Dibenzo(a,h)anthracene	26.532	278	3698395	39.481	ng	98
92) Benzo(g,h,i)perylene	27.291	276	3568421	38.936	ng	98

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

Quant Time: Mar 21 04:03:58 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM030823.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Mar 17 05:25:34 2023
Response via : Initial Calibration

