

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM033023\
 Data File : BM039234.D
 Acq On : 30 Mar 2023 19:46
 Operator : CG/JU
 Sample : 02023-01MSD
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 TP-11-TEST-PIT-ON-CENTRALMSD

Quant Time: Mar 31 02:41:00 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM032923.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Mar 30 05:15:45 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.922	152	289123	20.000	ng	0.00
21) Naphthalene-d8	10.733	136	1095097	20.000	ng	0.00
39) Acenaphthene-d10	14.568	164	585611	20.000	ng	0.00
64) Phenanthrene-d10	17.315	188	1069162	20.000	ng	0.00
76) Chrysene-d12	21.503	240	730357	20.000	ng	0.00
86) Perylene-d12	23.915	264	880994	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.487	112	1949826	101.604	ng	0.00
7) Phenol-d6	7.092	99	2405230	96.902	ng	0.00
23) Nitrobenzene-d5	9.092	82	1490563	66.154	ng	0.00
42) 2,4,6-Tribromophenol	16.062	330	671766	90.367	ng	0.00
45) 2-Fluorobiphenyl	13.192	172	2847554	65.883	ng	0.00
79) Terphenyl-d14	19.939	244	2916654	73.781	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	3.351	88	352670	44.961	ng	99
3) Pyridine	3.763	79	902343	41.063	ng	98
4) n-Nitrosodimethylamine	3.669	42	401822	46.560	ng	100
6) Aniline	7.251	93	935967	30.467	ng	100
8) 2-Chlorophenol	7.486	128	927864	47.240	ng	99
9) Benzaldehyde	7.057	77	600001	42.446	ng	98
10) Phenol	7.122	94	1160747	46.664	ng	99
11) bis(2-Chloroethyl)ether	7.345	93	939055	46.927	ng	100
12) 1,3-Dichlorobenzene	7.810	146	1011521	48.560	ng	99
13) 1,4-Dichlorobenzene	7.957	146	1021941	48.520	ng	98
14) 1,2-Dichlorobenzene	8.275	146	969494	48.138	ng	100
15) Benzyl Alcohol	8.169	79	782039	45.808	ng	99
16) 2,2'-oxybis(1-Chloropr...	8.463	45	1199154	46.850	ng	100
17) 2-Methylphenol	8.375	107	737385	42.371	ng	99
18) Hexachloroethane	9.004	117	400787	49.192	ng	97
19) n-Nitroso-di-n-propyla...	8.739	70	685693	43.215	ng	99
20) 3+4-Methylphenols	8.704	107	987369	42.222	ng	99
22) Acetophenone	8.757	105	1363087	46.553	ng	# 99
24) Nitrobenzene	9.139	77	1007487	44.954	ng	99
25) Isophorone	9.663	82	1826188	43.445	ng	100
26) 2-Nitrophenol	9.845	139	483246	46.646	ng	98
27) 2,4-Dimethylphenol	9.910	122	800377	46.498	ng	99
28) bis(2-Chloroethoxy)met...	10.151	93	1156270	45.442	ng	100
29) 2,4-Dichlorophenol	10.386	162	772798	46.015	ng	99
30) 1,2,4-Trichlorobenzene	10.598	180	844270	46.989	ng	98
31) Naphthalene	10.786	128	2672546	45.291	ng	100
32) Benzoic acid	10.063	122	586592	43.723	ng	98
33) 4-Chloroaniline	10.904	127	336689	13.024	ng	99
34) Hexachlorobutadiene	11.069	225	499984	47.324	ng	100
35) Caprolactam	11.698	113	225383	38.650	ng	96
36) 4-Chloro-3-methylphenol	12.027	107	780177	42.104	ng	98
37) 2-Methylnaphthalene	12.392	142	1751308	42.634	ng	99
38) 1-Methylnaphthalene	12.616	142	1623095	42.198	ng	100
40) 1,2,4,5-Tetrachloroben...	12.763	216	874076	49.023	ng	100
41) Hexachlorocyclopentadiene	12.739	237	718072	73.719	ng	97
43) 2,4,6-Trichlorophenol	13.004	196	563756	46.363	ng	99

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM033023\
 Data File : BM039234.D
 Acq On : 30 Mar 2023 19:46
 Operator : CG/JU
 Sample : 02023-01MSD
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 TP-11-TEST-PIT-ON-CENTRALMSD

Quant Time: Mar 31 02:41:00 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM032923.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Mar 30 05:15:45 2023
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	13.074	196	598576	41.991	ng	99
46) 1,1'-Biphenyl	13.404	154	2186506	47.282	ng	99
47) 2-Chloronaphthalene	13.445	162	1672810	47.774	ng	99
48) 2-Nitroaniline	13.657	65	502743	42.666	ng	97
49) Acenaphthylene	14.292	152	2678611	46.545	ng	99
50) Dimethylphthalate	14.033	163	1921706	42.580	ng	99
51) 2,6-Dinitrotoluene	14.151	165	417362	43.004	ng	98
52) Acenaphthene	14.633	154	1580059	46.498	ng	100
53) 3-Nitroaniline	14.480	138	291872	26.014	ng	97
54) 2,4-Dinitrophenol	14.692	184	482715	83.986	ng	96
55) Dibenzofuran	14.968	168	2402479	43.914	ng	99
56) 4-Nitrophenol	14.792	139	761773	88.048	ng	95
57) 2,4-Dinitrotoluene	14.939	165	558666	42.552	ng	94
58) Fluorene	15.615	166	1990036	45.937	ng	99
59) 2,3,4,6-Tetrachlorophenol	15.198	232	494852	43.637	ng	# 100
60) Diethylphthalate	15.392	149	1941202	42.733	ng	100
61) 4-Chlorophenyl-phenyle...	15.609	204	935328	43.565	ng	99
62) 4-Nitroaniline	15.645	138	418395	36.339	ng	99
63) Azobenzene	15.904	77	2269158	49.859	ng	100
65) 4,6-Dinitro-2-methylph...	15.704	198	302260	42.973	ng	93
66) n-Nitrosodiphenylamine	15.827	169	1588967	47.163	ng	99
67) 4-Bromophenyl-phenylether	16.504	248	537928	47.112	ng	98
68) Hexachlorobenzene	16.621	284	597342	48.249	ng	98
69) Atrazine	16.780	200	495966	46.479	ng	99
70) Pentachlorophenol	16.968	266	736041	82.865	ng	100
71) Phenanthrene	17.362	178	3022520	51.617	ng	100
72) Anthracene	17.451	178	2828841	47.378	ng	99
73) Carbazole	17.721	167	2370919	42.495	ng	100
74) Di-n-butylphthalate	18.286	149	3116817	45.161	ng	99
75) Fluoranthene	19.380	202	3181420	47.662	ng	99
77) Benzidine	19.562	184	727443	37.761	ng	99
78) Pyrene	19.739	202	3691124	70.992	ng	100
80) Butylbenzylphthalate	20.633	149	1198324	50.308	ng	96
81) Benzo(a)anthracene	21.486	228	2668394	51.804	ng	99
82) 3,3'-Dichlorobenzidine	21.421	252	685086	39.926	ng	99
83) Chrysene	21.538	228	2476683	49.807	ng	100
84) Bis(2-ethylhexyl)phtha...	21.409	149	1761691	50.077	ng	100
85) Di-n-octyl phthalate	22.338	149	2825192	44.574	ng	100
87) Indeno(1,2,3-cd)pyrene	26.432	276	3495463	53.850	ng	99
88) Benzo(b)fluoranthene	23.180	252	2697840	50.047	ng	100
89) Benzo(k)fluoranthene	23.233	252	2589964	47.168	ng	99
90) Benzo(a)pyrene	23.809	252	2611241	49.310	ng	100
91) Dibenzo(a,h)anthracene	26.450	278	2777685	50.962	ng	100
92) Benzo(g,h,i)perylene	27.203	276	2948067	54.999	ng	99

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

Instrument :
BNA_M
ClientSampleId :
TP-11-TEST-PIT-ON-CENTRALMSD

[illegible]