

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM040324\
 Data File : BM044892.D
 Acq On : 03 Apr 2024 13:02
 Operator : MA/JU
 Sample : P1908-03DL 5X
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 PT-BN-WPDL

Quant Time: Apr 03 13:35:13 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM040224.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Apr 03 00:28:57 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.645	152	71416	20.000	ng	0.00
21) Naphthalene-d8	10.422	136	289506	20.000	ng	0.00
39) Acenaphthene-d10	14.298	164	187792	20.000	ng	0.00
64) Phenanthrene-d10	17.057	188	404911	20.000	ng	0.00
76) Chrysene-d12	21.262	240	425648	20.000	ng	0.00
86) Perylene-d12	23.497	264	549672	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.257	112	140379	29.677	ng	0.00
7) Phenol-d6	6.828	99	175445	28.098	ng	0.00
23) Nitrobenzene-d5	8.804	82	108416	18.442	ng	-0.01
42) 2,4,6-Tribromophenol	15.798	330	65414	23.837	ng	0.00
45) 2-Fluorobiphenyl	12.910	172	263879	19.058	ng	0.00
79) Terphenyl-d14	19.703	244	481180	19.316	ng	0.00
Target Compounds						Qvalue
11) bis(2-Chloroethyl)ether	7.092	93	58944	11.698	ng	95
18) Hexachloroethane	8.698	117	32830	15.808	ng	98
19) n-Nitroso-di-n-propyla...	8.457	70	71192	17.084	ng	96
24) Nitrobenzene	8.851	77	125317	21.797	ng	98
28) bis(2-Chloroethoxy)met...	9.857	93	152314	23.590	ng	95
31) Naphthalene	10.475	128	191179	12.218	ng	99
34) Hexachlorobutadiene	10.739	225	54965	20.654	ng	95
37) 2-Methylnaphthalene	12.098	142	146206	13.004	ng	98
38) 1-Methylnaphthalene	12.316	142	76543	7.191	ng	97
47) 2-Chloronaphthalene	13.163	162	377667	35.373	ng	100
49) Acenaphthylene	14.016	152	290165	16.330	ng	99
51) 2,6-Dinitrotoluene	13.892	165	87983	28.275	ng	98
52) Acenaphthene	14.363	154	73584	6.870	ng	96
55) Dibenzofuran	14.698	168	217169	12.299	ng	96
57) 2,4-Dinitrotoluene	14.692	165	100913	23.189	ng	95
58) Fluorene	15.351	166	362879	25.358	ng	100
60) Diethylphthalate	15.139	149	324244	21.286	ng	99
62) 4-Nitroaniline	15.398	138	67914	20.906	ng	98
66) n-Nitrosodiphenylamine	15.574	169	155132	12.662	ng	98
68) Hexachlorobenzene	16.351	284	67247	13.417	ng	98
71) Phenanthrene	17.098	178	355825	15.810	ng	99
72) Anthracene	17.186	178	479265	21.031	ng	100
74) Di-n-butylphthalate	18.039	149	913007	33.058	ng	99
75) Fluoranthene	19.127	202	926302	34.779	ng	99
78) Pyrene	19.492	202	507852	17.918	ng	99
80) Butylbenzylphthalate	20.415	149	354013	27.067	ng	98
83) Chrysene	21.303	228	319814	11.474	ng	99
84) Bis(2-ethylhexyl)phtha...	21.197	149	455820	21.818	ng	99
87) Indeno(1,2,3-cd)pyrene	25.750	276	1239073	30.043	ng	95
88) Benzo(b)fluoranthene	22.833	252	996202	30.566	ng	99
89) Benzo(k)fluoranthene	22.874	252	247258	7.639	ng	99
90) Benzo(a)pyrene	23.403	252	459727	14.563	ng	99
91) Dibenzo(a,h)anthracene	25.762	278	324647	9.348	ng	98
92) Benzo(g,h,i)perylene	26.426	276	194529	5.831	ng	98

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

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