

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM040324\
 Data File : BM044893.D
 Acq On : 03 Apr 2024 15:03
 Operator : MA/JU
 Sample : P1908-06DL 5X
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 PT-ACIDS-WPDL

Quant Time: Apr 03 15:46:35 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	78837	20.000	ng	0.00
21) Naphthalene-d8	10.422	136	315273	20.000	ng	0.00
39) Acenaphthene-d10	14.298	164	201881	20.000	ng	0.00
64) Phenanthrene-d10	17.057	188	433973	20.000	ng	0.00
76) Chrysene-d12	21.268	240	442414	20.000	ng	0.00
86) Perylene-d12	23.503	264	586953	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.257	112	143487	27.479	ng	0.00
7) Phenol-d6	6.822	99	178694	25.924	ng	-0.01
23) Nitrobenzene-d5	8.804	82	108982	17.024	ng	-0.01
42) 2,4,6-Tribromophenol	15.798	330	74337	25.198	ng	0.00
45) 2-Fluorobiphenyl	12.910	172	261868	17.593	ng	0.00
79) Terphenyl-d14	19.703	244	456766	17.641	ng	0.00
Target Compounds						Qvalue
8) 2-Chlorophenol	7.210	128	166385	29.839	ng	98
10) Phenol	6.851	94	174643	25.687	ng	98
17) 2-Methylphenol	8.092	107	59489	12.019	ng	100
20) 3+4-Methylphenols	8.410	107	192845	28.034	ng	91
26) 2-Nitrophenol	9.551	139	84232	29.370	ng	98
27) 2,4-Dimethylphenol	9.616	122	39625	8.266	ng	96
29) 2,4-Dichlorophenol	10.086	162	34513	7.074	ng	98
36) 4-Chloro-3-methylphenol	11.727	107	108378	20.505	ng	96
43) 2,4,6-Trichlorophenol	12.722	196	61231	15.608	ng	95
54) 2,4-Dinitrophenol	14.445	184	44389	23.572	ng	96
56) 4-Nitrophenol	14.539	139	84241	27.724	ng	98
59) 2,3,4,6-Tetrachlorophenol	14.933	232	37807	9.676	ng	99
65) 4,6-Dinitro-2-methylph...	15.451	198	50325	18.303	ng	97
70) Pentachlorophenol	16.704	266	76532	19.679	ng	96

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

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