

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM032122\
 Data File : BM034287.D
 Acq On : 21 Mar 2022 13:37
 Operator : CG/JU
 Sample : N1870-06DL 5X
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 PT-ACIDS-WPDL

Quant Time: Mar 21 17:55:01 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM031522.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Mar 16 02:38:26 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.942	152	135757	20.000	ng	0.00
21) Naphthalene-d8	10.754	136	564154	20.000	ng	-0.01
39) Acenaphthene-d10	14.583	164	369651	20.000	ng	0.00
64) Phenanthrene-d10	17.324	188	769454	20.000	ng	0.00
76) Chrysene-d12	21.494	240	765266	20.000	ng	0.00
86) Perylene-d12	23.906	264	748916	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.490	112	229574	30.337	ng	-0.01
7) Phenol-d6	7.095	99	301591	27.835	ng	-0.01
23) Nitrobenzene-d5	9.101	82	209051	18.041	ng	-0.01
42) 2,4,6-Tribromophenol	16.065	330	137959	27.689	ng	0.00
45) 2-Fluorobiphenyl	13.213	172	524117	19.375	ng	0.00
79) Terphenyl-d14	19.936	244	865859	19.660	ng	-0.01
Target Compounds						
						Qvalue
8) 2-Chlorophenol	7.501	128	59719	6.747	ng	97
10) Phenol	7.119	94	269604	25.019	ng	98
20) 3+4-Methylphenols	8.707	107	134599	12.895	ng	91
26) 2-Nitrophenol	9.866	139	131080	27.322	ng	99
27) 2,4-Dimethylphenol	9.925	122	140546	18.709	ng	98
29) 2,4-Dichlorophenol	10.395	162	219405	25.137	ng	100
36) 4-Chloro-3-methylphenol	12.030	107	315468	31.539	ng	97
43) 2,4,6-Trichlorophenol	13.019	196	116088	15.073	ng	99
44) 2,4,5-Trichlorophenol	13.083	196	274707	33.231	ng	99
54) 2,4-Dinitrophenol	14.683	184	73305	22.673	ng	96
56) 4-Nitrophenol	14.777	139	58543	12.473	ng	97
65) 4,6-Dinitro-2-methylph...	15.695	198	32142	6.424	ng	95
70) Pentachlorophenol	16.971	266	171714	27.216	ng	99

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

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