

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF091624\
 Data File : BF139380.D
 Acq On : 16 Sep 2024 13:13
 Operator : RC/JU
 Sample : P3845-06DL 5X
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PT-ACIDS-WPDL

Quant Time: Sep 16 14:01:35 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF091324.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Sep 14 02:58:23 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.875	152	53717	20.000	ng	0.00
21) Naphthalene-d8	8.157	136	197693	20.000	ng	0.00
39) Acenaphthene-d10	9.910	164	111417	20.000	ng	0.00
64) Phenanthrene-d10	11.392	188	205713	20.000	ng	0.00
76) Chrysene-d12	14.027	240	127464	20.000	ng	0.00
86) Perylene-d12	15.498	264	107390	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.487	112	81036	24.610	ng	0.00
7) Phenol-d6	6.492	99	107117	24.784	ng	-0.01
23) Nitrobenzene-d5	7.433	82	72704	17.289	ng	-0.01
42) 2,4,6-Tribromophenol	10.692	330	30178	25.442	ng	0.00
45) 2-Fluorobiphenyl	9.227	172	144091	18.138	ng	0.00
79) Terphenyl-d14	12.974	244	147214	18.957	ng	0.00
Target Compounds						
						Qvalue
8) 2-Chlorophenol	6.657	128	97582	28.789	ng	99
10) Phenol	6.510	94	100447	22.819	ng	98
17) 2-Methylphenol	7.122	107	53975	20.043	ng	99
26) 2-Nitrophenol	7.769	139	33311	19.755	ng	98
27) 2,4-Dimethylphenol	7.804	122	28375	12.757	ng	98
29) 2,4-Dichlorophenol	8.010	162	27945	9.890	ng	98
36) 4-Chloro-3-methylphenol	8.698	107	92018	28.495	ng	99
43) 2,4,6-Trichlorophenol	9.145	196	62186	28.947	ng	98
44) 2,4,5-Trichlorophenol	9.180	196	26419	11.658	ng	98
54) 2,4-Dinitrophenol	9.963	184	19289	21.368	ng	94
56) 4-Nitrophenol	10.010	139	38922	29.054	ng	95
59) 2,3,4,6-Tetrachlorophenol	10.227	232	37016	19.746	ng	98
65) 4,6-Dinitro-2-methylph...	10.492	198	21982	20.246	ng	98
70) Pentachlorophenol	11.192	266	9184	6.397	ng	97

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

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