

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110420\
 Data File : BP003806.D
 Acq On : 04 Nov 2020 15:15
 Operator : CG/JU
 Sample : L3971-06DL 5X
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 PT-ACIDS-WPDL

Quant Time: Nov 05 14:55:45 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-BP110320.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Nov 05 14:31:45 2020
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.293	152	171269	20.00	ng	0.00
21) Naphthalene-d8	11.128	136	661987	20.00	ng	# 0.00
39) Acenaphthene-d10	14.910	164	401411	20.00	ng	0.00
64) Phenanthrene-d10	17.634	188	835817	20.00	ng	# 0.00
76) Chrysene-d12	21.651	240	821387	20.00	ng	0.00
86) Perylene-d12	24.122	264	833620	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.811	112	246891	24.06	ng	0.00
7) Phenol-d6	7.452	99	321344	21.81	ng	0.00
23) Nitrobenzene-d5	9.458	82	205664	15.22	ng	0.00
42) 2,4,6-Tribromophenol	16.387	330	108880	21.69	ng	0.00
45) 2-Fluorobiphenyl	13.540	172	487394	16.97	ng	0.00
79) Terphenyl-d14	20.122	244	678889	15.97	ng	0.00
Target Compounds						Qvalue
8) 2-Chlorophenol	7.864	128	194441	17.85	ng	# 81
10) Phenol	7.481	94	367296	25.84	ng	84
17) 2-Methylphenol	8.746	107	86564	9.27	ng	# 92
20) 3+4-Methylphenols	9.069	107	189189	15.19	ng	92
26) 2-Nitrophenol	10.228	139	93153	17.76	ng	# 74
27) 2,4-Dimethylphenol	10.287	122	193512	22.12	ng	92
29) 2,4-Dichlorophenol	10.775	162	250433	23.75	ng	96
36) 4-Chloro-3-methylphenol	12.381	107	84594	7.43	ng	95
43) 2,4,6-Trichlorophenol	13.363	196	198038	24.11	ng	100
44) 2,4,5-Trichlorophenol	13.434	196	55048	6.37	ng	97
54) 2,4-Dinitrophenol	14.987	184	45631	20.29	ng	# 82
65) 4,6-Dinitro-2-methylph...	16.010	198	87518	23.22	ng	# 79
70) Pentachlorophenol	17.304	266	39858	7.47	ng	99

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

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