

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM100421\
 Data File : BM032322.D
 Acq On : 04 Oct 2021 14:20
 Operator : CG/JU
 Sample : M3702-06DL 5X
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 PT-ACIDS-WPDL

Quant Time: Oct 04 14:50:05 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM100221.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Oct 01 18:34:58 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.637	152	329868	20.00	ng	0.00
21) Naphthalene-d8	10.425	136	1294720	20.00	ng	# 0.00
39) Acenaphthene-d10	14.283	164	789297	20.00	ng	0.00
64) Phenanthrene-d10	17.007	188	1629321	20.00	ng	# 0.00
76) Chrysene-d12	21.177	240	1479036	20.00	ng	0.00
86) Perylene-d12	23.330	264	1381939	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.207	112	598572	30.57	ng	0.00
7) Phenol-d6	6.819	99	752369	28.10	ng	-0.01
23) Nitrobenzene-d5	8.784	82	425059	18.26	ng	-0.02
42) 2,4,6-Tribromophenol	15.765	330	244336	27.65	ng	0.00
45) 2-Fluorobiphenyl	12.901	172	1065426	20.07	ng	0.00
79) Terphenyl-d14	19.624	244	1539814	21.93	ng	0.00
Target Compounds						
						Qvalue
8) 2-Chlorophenol	7.207	128	631023	29.57	ng	91
10) Phenol	6.848	94	722342	28.02	ng	82
17) 2-Methylphenol	8.095	107	282302	16.21	ng	# 90
20) 3+4-Methylphenols	8.413	107	854021	36.34	ng	86
26) 2-Nitrophenol	9.542	139	250133	23.94	ng	# 75
27) 2,4-Dimethylphenol	9.619	122	547204	31.05	ng	93
29) 2,4-Dichlorophenol	10.083	162	365993	18.65	ng	95
36) 4-Chloro-3-methylphenol	11.719	107	338315	16.01	ng	96
43) 2,4,6-Trichlorophenol	12.713	196	196726	12.51	ng	95
44) 2,4,5-Trichlorophenol	12.783	196	213982	12.65	ng	# 91
56) 4-Nitrophenol	14.495	139	230457	22.03	ng	# 71
65) 4,6-Dinitro-2-methylph...	15.401	198	97811	10.67	ng	# 78
70) Pentachlorophenol	16.677	266	160654	13.94	ng	98

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

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