

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM100421\
 Data File : BM032321.D
 Acq On : 04 Oct 2021 13:28
 Operator : CG/JU
 Sample : M3702-06
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 PT-ACIDS-WP

Quant Time: Oct 04 13:58:35 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	310021	20.00	ng	0.00
21) Naphthalene-d8	10.425	136	1226737	20.00	ng	# 0.00
39) Acenaphthene-d10	14.283	164	730043	20.00	ng	0.00
64) Phenanthrene-d10	17.012	188	1500876	20.00	ng	0.00
76) Chrysene-d12	21.177	240	1325406	20.00	ng	0.00
86) Perylene-d12	23.330	264	1220964	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.207	112	2830919	153.86	ng	0.00
7) Phenol-d6	6.831	99	3599611	143.06	ng	0.00
23) Nitrobenzene-d5	8.795	82	2194911	99.51	ng	0.00
42) 2,4,6-Tribromophenol	15.771	330	1175440	143.82	ng	0.00
45) 2-Fluorobiphenyl	12.901	172	4591772	93.52	ng	0.00
79) Terphenyl-d14	19.630	244	5881788	93.46	ng	0.00
Target Compounds						
						Qvalue
8) 2-Chlorophenol	7.207	128	2954271	147.31	ng	93
10) Phenol	6.860	94	3349108	138.24	ng	83
17) 2-Methylphenol	8.095	107	1405092	85.82	ng	# 90
20) 3+4-Methylphenols	8.431	107	3912968	177.15	ng	86
26) 2-Nitrophenol	9.548	139	1388749	140.30	ng	# 74
27) 2,4-Dimethylphenol	9.625	122	2502778	149.88	ng	95
29) 2,4-Dichlorophenol	10.083	162	1827984	98.29	ng	97
36) 4-Chloro-3-methylphenol	11.725	107	1708145	85.31	ng	96
43) 2,4,6-Trichlorophenol	12.713	196	1018589	70.00	ng	96
44) 2,4,5-Trichlorophenol	12.783	196	1113038	71.16	ng	# 92
56) 4-Nitrophenol	14.513	139	1351774	139.69	ng	# 67
65) 4,6-Dinitro-2-methylph...	15.407	198	738023	87.36	ng	# 72
70) Pentachlorophenol	16.677	266	874366	82.38	ng	98

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

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