

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG031120\
 Data File : BG044724.D
 Acq On : 11 Mar 2020 17:52
 Operator : CG/JU
 Sample : L1798-06DL 5X
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PT-ACIDS-WPDL

Quant Time: Mar 12 03:46:55 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG031020.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Mar 10 16:27:39 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.06	152	87181	20.00	ng	0.00
21) Naphthalene-d8	10.87	136	357157	20.00	ng	0.00
39) Acenaphthene-d10	14.69	164	246952	20.00	ng	0.00
64) Phenanthrene-d10	17.44	188	562958	20.00	ng	0.00
76) Chrysene-d12	21.70	240	554790	20.00	ng	0.00
87) Perylene-d12	24.93	264	622934	20.00	ng	0.00

System Monitoring Compounds						
5) 2-Fluorophenol	5.63	112	172493	30.80	ng	0.00
7) Phenol-d6	7.21	99	256697	31.55	ng	0.00
23) Nitrobenzene-d5	9.21	82	158529	19.48	ng	0.00
42) 2,4,6-Tribromophenol	16.18	330	93589	28.94	ng	0.00
45) 2-Fluorobiphenyl	13.32	172	371313	21.53	ng	0.00
79) Terphenyl-d14	20.05	244	697663	23.52	ng	0.00

Target Compounds				Qvalue		
8) 2-Chlorophenol	7.63	128	160399	28.691	ng	94
17) 2-Methylphenol	8.50	107	90514	16.587	ng	94
20) 3+4-Methylphenols	8.81	107	213274	28.352	ng	91
26) 2-Nitrophenol	9.98	139	42359	18.303	ng	94
27) 2,4-Dimethylphenol	10.04	122	83132	16.070	ng	95
29) 2,4-Dichlorophenol	10.51	162	99153	16.410	ng	96
36) 4-Chloro-3-methylphenol	12.13	107	157946	21.918	ng	97
43) 2,4,6-Trichlorophenol	13.13	196	55638	10.308	ng	96
44) 2,4,5-Trichlorophenol	13.19	196	67117	11.991	ng	91
54) 2,4-Dinitrophenol	14.78	184	35696	23.760	ng	# 57
56) 4-Nitrophenol	14.88	139	87223	25.074	ng	99
65) 4,6-Dinitro-2-methylphenol	15.81	198	70839	30.163	ng	95
70) Pentachlorophenol	17.09	266	115432	27.477	ng	98

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

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