

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP021022\
 Data File : BP009106.D
 Acq On : 10 Feb 2022 18:56
 Operator : CG/JU
 Sample : N1357-01DL 5X
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 GROUNDWATERDL

Quant Time: Feb 11 02:26:35 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP020722.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Feb 07 15:13:17 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.799	152	250616	20.000	ng	0.00
21) Naphthalene-d8	10.599	136	964729	20.000	ng	0.00
39) Acenaphthene-d10	14.446	164	639178	20.000	ng	0.00
64) Phenanthrene-d10	17.204	188	1362924	20.000	ng	0.00
76) Chrysene-d12	21.298	240	1247225	20.000	ng	0.00
86) Perylene-d12	23.704	264	1362243	20.000	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.405	112	147545	10.502	ng	0.01
7) Phenol-d6	7.005	99	123457	6.669	ng	0.02
23) Nitrobenzene-d5	8.963	82	296013	17.304	ng	0.00
42) 2,4,6-Tribromophenol	15.945	330	227745	26.686	ng	0.00
45) 2-Fluorobiphenyl	13.069	172	793383	17.376	ng	0.00
79) Terphenyl-d14	19.769	244	1194439	17.222	ng	0.00
Target Compounds						
						Qvalue
6) Aniline	7.134	93	217609	10.356	ng	99
10) Phenol	7.028	94	396006	21.276	ng	94
15) Benzyl Alcohol	8.069	79	24715	2.097	ng	92
17) 2-Methylphenol	8.269	107	254112	20.364	ng	97
20) 3+4-Methylphenols	8.611	107	84911	4.972	ng	88
22) Acetophenone	8.634	105	49818	2.209	ng	# 83
27) 2,4-Dimethylphenol	9.805	122	327316	26.172	ng	98
31) Naphthalene	10.652	128	554001	11.259	ng	99
33) 4-Chloroaniline	10.858	127	36721960	1848.269	ng	95
60) Diethylphthalate	15.275	149	155447	3.608	ng	99
80) Butylbenzylphthalate	20.445	149	278038	9.114	ng	98

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

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