

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG101617\
 Data File : BG029279.D
 Acq On : 16 Oct 2017 18:46
 Operator : SJ/JU
 Sample : I5266-06DL 5X
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PT-ACIDS-WPDL

Quant Time: Oct 17 17:08:03 2017
 Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG101617.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Oct 16 17:32:15 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.17	152	63279	20.00	ng	0.00
21) Naphthalene-d8	10.99	136	302192	20.00	ng	0.00
38) Acenaphthene-d10	14.79	164	198360	20.00	ng	0.00
63) Phenanthrene-d10	17.53	188	452668	20.00	ng	0.00
75) Chrysene-d12	21.80	240	470257	20.00	ng	0.00
86) Perylene-d12	25.11	264	483209	20.00	ng	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
5) 2-Fluorophenol	5.72	112	102741	27.12	ng	0.00
7) Phenol-d6	7.32	99	139376	26.21	ng	0.00
23) Nitrobenzene-d5	9.33	82	79642	16.75	ng	0.00
41) 2,4,6-Tribromophenol	16.27	330	62373	24.35	ng	0.00
44) 2-Fluorobiphenyl	13.41	172	240341	17.77	ng	0.00
78) Terphenyl-d14	20.12	244	334890	16.20	ng	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
8) 2-Chlorophenol	7.73	128	162833	37.503	ng	95
10) Phenol	7.35	94	200937	35.054	ng	91
17) 2-Methylphenol	8.61	107	64346	16.898	ng	95
20) 3+4-Methylphenols	8.93	107	98520	18.362	ng	91
26) 2-Nitrophenol	10.09	139	42909	16.791	ng	91
27) 2,4-Dimethylphenol	10.15	122	82982	17.948	ng	90
29) 2,4-Dichlorophenol	10.63	162	151379	30.703	ng	93
36) 4-Chloro-3-methylphenol	12.24	107	47527	8.764	ng	87
42) 2,4,6-Trichlorophenol	13.23	196	76044	16.727	ng	100
43) 2,4,5-Trichlorophenol	13.30	196	58469	12.523	ng	95
55) 4-Nitrophenol	14.98	139	88773	30.757	ng	# 81
64) 4,6-Dinitro-2-methylphenol	15.90	198	51045	22.764	ng	89
69) Pentachlorophenol	17.18	266	74387	22.009	ng	99

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

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