

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072823\
 Data File : BP016438.D
 Acq On : 28 Jul 2023 23:42
 Operator : MA/JU
 Sample : 03813-08
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 PLAUDERVILLE-3

Quant Time: Jul 29 03:48:14 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP072723.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jul 28 01:48:25 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.087	152	371177	20.000	ng	0.00
21) Naphthalene-d8	10.922	136	1565531	20.000	ng	0.00
39) Acenaphthene-d10	14.728	164	921423	20.000	ng	0.00
64) Phenanthrene-d10	17.492	188	2062453	20.000	ng	0.00
76) Chrysene-d12	21.580	240	1744133	20.000	ng	#-0.01
86) Perylene-d12	24.174	264	1731039	20.000	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.605	112	2794085	122.807	ng	0.00
7) Phenol-d6	7.222	99	3476982	111.494	ng	0.00
23) Nitrobenzene-d5	9.281	82	2329782	83.306	ng	-0.01
42) 2,4,6-Tribromophenol	16.222	330	1386755	102.302	ng	0.00
45) 2-Fluorobiphenyl	13.357	172	5197029	80.862	ng	0.00
79) Terphenyl-d14	20.033	244	8452625	102.375	ng	-0.01
Target Compounds						
32) Benzoic acid	10.116	122	32635	7.997	ng	Qvalue 92

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

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