

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072823\
 Data File : BP016441.D
 Acq On : 29 Jul 2023 01:25
 Operator : MA/JU
 Sample : 03770-08
 Misc :
 ALS Vial : 29 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 HOWE-1-LP

Quant Time: Jul 29 03:49:21 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	334245	20.000	ng	0.00
21) Naphthalene-d8	10.922	136	1419886	20.000	ng	0.00
39) Acenaphthene-d10	14.728	164	869961	20.000	ng	0.00
64) Phenanthrene-d10	17.492	188	2053261	20.000	ng	0.00
76) Chrysene-d12	21.580	240	1858911	20.000	ng	#-0.01
86) Perylene-d12	24.174	264	1886137	20.000	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.605	112	2878567	140.500	ng	0.00
7) Phenol-d6	7.222	99	3583482	127.606	ng	0.00
23) Nitrobenzene-d5	9.281	82	2387902	94.142	ng	-0.01
42) 2,4,6-Tribromophenol	16.222	330	1460994	114.154	ng	0.00
45) 2-Fluorobiphenyl	13.357	172	5384822	88.740	ng	0.00
79) Terphenyl-d14	20.033	244	9154696	104.032	ng	-0.01
Target Compounds						
32) Benzoic acid	10.110	122	19620	7.420	ng	Qvalue 86

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

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