

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072823\
 Data File : BP016436.D
 Acq On : 28 Jul 2023 22:33
 Operator : MA/JU
 Sample : 03797-05
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 HANOVER-ST-2

Quant Time: Jul 29 00:44:33 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	310679	20.000	ng	0.00
21) Naphthalene-d8	10.922	136	1301427	20.000	ng	0.00
39) Acenaphthene-d10	14.728	164	780757	20.000	ng	0.00
64) Phenanthrene-d10	17.492	188	1815695	20.000	ng	0.00
76) Chrysene-d12	21.580	240	1675213	20.000	ng	-0.01
86) Perylene-d12	24.174	264	1740041	20.000	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.599	112	2097520	110.143	ng	0.00
7) Phenol-d6	7.222	99	2648311	101.458	ng	0.00
23) Nitrobenzene-d5	9.281	82	1718612	73.923	ng	-0.01
42) 2,4,6-Tribromophenol	16.222	330	1153384	100.416	ng	0.00
45) 2-Fluorobiphenyl	13.357	172	3931368	72.190	ng	0.00
79) Terphenyl-d14	20.034	244	7692213	96.998	ng	-0.01
Target Compounds						
32) Benzoic acid	10.116	122	40798	8.860	ng	Qvalue 92

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

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