

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP061324\  
 Data File : BP020614.D  
 Acq On : 13 Jun 2024 16:38  
 Operator : MA/JU  
 Sample : P2857-02  
 Misc :  
 ALS Vial : 6 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 ETGI-355

Quant Time: Jun 13 17:08:03 2024  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\8270E-BP052924.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Thu May 30 03:41:03 2024  
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.752	152	211187	20.000	ng	-0.02
21) Naphthalene-d8	10.516	136	814100	20.000	ng	-0.02
39) Acenaphthene-d10	14.375	164	534291	20.000	ng	-0.02
64) Phenanthrene-d10	17.181	188	1137478	20.000	ng	-0.02
76) Chrysene-d12	21.633	240	1063693	20.000	ng	-0.04
86) Perylene-d12	24.986	264	1211332	20.000	ng	-0.04
System Monitoring Compounds						
5) 2-Fluorophenol	5.369	112	1045907	88.672	ng	0.00
7) Phenol-d6	6.928	99	1524454	91.646	ng	0.00
23) Nitrobenzene-d5	8.899	82	1056250	71.022	ng	-0.02
42) 2,4,6-Tribromophenol	15.886	330	467521	84.315	ng	-0.02
45) 2-Fluorobiphenyl	12.987	172	2567808	71.644	ng	-0.02
79) Terphenyl-d14	19.916	244	4385849	68.640	ng	-0.03

Target Compounds Qvalue

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

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