

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP030323\
 Data File : BP013945.D
 Acq On : 03 Mar 2023 15:24
 Operator : CG/JU
 Sample : 01763-01
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 CM-A-CRUSHED-MASONRY-PILE-1

Quant Time: Mar 04 00:56:36 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP021623.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Mar 02 04:26:35 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.093	152	105983	20.000	ng	0.00
21) Naphthalene-d8	10.916	136	368835	20.000	ng	0.00
39) Acenaphthene-d10	14.728	164	217397	20.000	ng	0.00
64) Phenanthrene-d10	17.481	188	487657	20.000	ng	0.00
76) Chrysene-d12	21.557	240	564217	20.000	ng	-0.01
86) Perylene-d12	24.104	264	668684	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.628	112	521891	81.158	ng	0.00
7) Phenol-d6	7.240	99	724793	88.005	ng	-0.01
23) Nitrobenzene-d5	9.263	82	493251	72.215	ng	0.00
42) 2,4,6-Tribromophenol	16.216	330	199050	62.899	ng	-0.01
45) 2-Fluorobiphenyl	13.352	172	1070879	64.700	ng	0.00
79) Terphenyl-d14	20.016	244	1968763	67.135	ng	0.00

Target Compounds	Qvalue

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

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