

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM102622\
 Data File : BM037250.D
 Acq On : 26 Oct 2022 22:49
 Operator : CG/JU
 Sample : N5218-11
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SP-2

Quant Time: Oct 27 07:52:55 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM102622.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Oct 27 04:49:27 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.887	152	27912	20.000	ng	0.00
21) Naphthalene-d8	10.687	136	138676	20.000	ng	0.00
39) Acenaphthene-d10	14.522	164	120168	20.000	ng	0.00
64) Phenanthrene-d10	17.263	188	405923	20.000	ng	0.00
76) Chrysene-d12	21.433	240	902945	20.000	ng	0.00
86) Perylene-d12	23.803	264	1291836	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.457	112	194843	143.118	ng	0.00
7) Phenol-d6	7.063	99	290202	154.014	ng	0.00
23) Nitrobenzene-d5	9.051	82	134985	55.569	ng	0.00
42) 2,4,6-Tribromophenol	16.010	330	550471	369.535	ng	0.00
45) 2-Fluorobiphenyl	13.145	172	413717	48.621	ng	0.00
79) Terphenyl-d14	19.880	244	1918111	43.456	ng	0.00

Target Compounds	Qvalue

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

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