

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG090222\
 Data File : BG054840.D
 Acq On : 2 Sep 2022 14:49
 Operator : CG/JU
 Sample : N4466-01
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SASP2-T-5-(10-17)

Quant Time: Sep 03 00:27:03 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG082522.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Aug 25 17:50:55 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.148	152	30564	20.000	ng	0.00
21) Naphthalene-d8	10.974	136	123515	20.000	ng	0.00
39) Acenaphthene-d10	14.781	164	88570	20.000	ng	0.00
64) Phenanthrene-d10	17.531	188	229397	20.000	ng	0.00
76) Chrysene-d12	21.820	240	229491	20.000	ng	0.00
86) Perylene-d12	25.192	264	235739	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.686	112	185757	105.834	ng	0.00
7) Phenol-d6	7.296	99	254709	106.113	ng	0.00
23) Nitrobenzene-d5	9.323	82	147175	62.552	ng	0.00
42) 2,4,6-Tribromophenol	16.268	330	128009	115.663	ng	0.00
45) 2-Fluorobiphenyl	13.406	172	371385	63.023	ng	0.00
79) Terphenyl-d14	20.128	244	712560	61.521	ng	0.00

Target Compounds	Qvalue

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

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