

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM110923\
 Data File : BM042677.D
 Acq On : 09 Nov 2023 18:34
 Operator : MA/JU
 Sample : 05267-01
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 ETGI-320

Quant Time: Nov 09 23:53:12 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM103023.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Nov 09 13:21:00 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.845	152	28853	20.000	ng	0.00
21) Naphthalene-d8	10.663	136	98993	20.000	ng	# 0.00
39) Acenaphthene-d10	14.498	164	57309	20.000	ng	0.00
64) Phenanthrene-d10	17.256	188	112395	20.000	ng	# 0.00
76) Chrysene-d12	21.439	240	106096	20.000	ng	# 0.00
86) Perylene-d12	23.809	264	150817	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.375	112	161265	90.203	ng	0.00
7) Phenol-d6	7.010	99	251433	102.546	ng	0.00
23) Nitrobenzene-d5	9.045	82	196893	92.718	ng	0.00
42) 2,4,6-Tribromophenol	15.998	330	30971	41.393	ng	0.00
45) 2-Fluorobiphenyl	13.121	172	374093	87.836	ng	0.00
79) Terphenyl-d14	19.874	244	523442	83.656	ng	0.00

Target Compounds	Qvalue

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

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