

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM111023\
 Data File : BM042687.D
 Acq On : 10 Nov 2023 12:34
 Operator : MA/JU
 Sample : 05279-04
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 TP-1

Quant Time: Nov 10 13:06:17 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	47656	20.000	ng	0.00
21) Naphthalene-d8	10.663	136	174876	20.000	ng	0.00
39) Acenaphthene-d10	14.504	164	120563	20.000	ng	0.00
64) Phenanthrene-d10	17.257	188	287697	20.000	ng	# 0.00
76) Chrysene-d12	21.439	240	233638	20.000	ng	# 0.00
86) Perylene-d12	23.815	264	92418	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.381	112	423962	143.576	ng	0.00
7) Phenol-d6	7.016	99	509446	125.797	ng	0.00
23) Nitrobenzene-d5	9.051	82	433878	115.658	ng	0.00
42) 2,4,6-Tribromophenol	16.004	330	340200	216.130	ng	0.00
45) 2-Fluorobiphenyl	13.122	172	994911	111.042	ng	0.00
79) Terphenyl-d14	19.874	244	1969046	142.902	ng	0.00

Target Compounds	Qvalue

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

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