

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM080223\
 Data File : BM041239.D
 Acq On : 02 Aug 2023 11:34
 Operator : MA/JU
 Sample : 03810-02
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 B-P3B

Quant Time: Aug 03 03:05:52 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM072823.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Aug 02 19:20:51 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.795	152	146183	20.000	ng	0.00
21) Naphthalene-d8	10.606	136	659664	20.000	ng	0.00
39) Acenaphthene-d10	14.465	164	452654	20.000	ng	0.00
64) Phenanthrene-d10	17.223	188	1042363	20.000	ng	0.00
76) Chrysene-d12	21.429	240	766472	20.000	ng	0.00
86) Perylene-d12	23.805	264	783883	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.371	112	821037	83.943	ng	0.00
7) Phenol-d6	6.983	99	1115639	87.996	ng	0.00
23) Nitrobenzene-d5	8.971	82	669523	47.200	ng	0.00
42) 2,4,6-Tribromophenol	15.970	330	554204	86.560	ng	0.00
45) 2-Fluorobiphenyl	13.082	172	1869236	52.166	ng	0.00
79) Terphenyl-d14	19.864	244	3340906	78.810	ng	0.00

Target Compounds	Qvalue

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

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