

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM082123\
 Data File : BM041447.D
 Acq On : 21 Aug 2023 14:27
 Operator : MA/JU
 Sample : 04082-02
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 277

Quant Time: Aug 22 03:36:55 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM081623.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Aug 22 03:32:15 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.757	152	157038	20.000	ng	0.00
21) Naphthalene-d8	10.557	136	627818	20.000	ng	0.00
39) Acenaphthene-d10	14.416	164	415131	20.000	ng	0.00
64) Phenanthrene-d10	17.180	188	956410	20.000	ng	0.00
76) Chrysene-d12	21.380	240	1011305	20.000	ng	0.00
86) Perylene-d12	23.727	264	1181821	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.346	112	1010563	107.053	ng	0.00
7) Phenol-d6	6.957	99	1173716	91.567	ng	0.00
23) Nitrobenzene-d5	8.939	82	913858	69.485	ng	0.00
42) 2,4,6-Tribromophenol	15.921	330	678711	103.024	ng	0.00
45) 2-Fluorobiphenyl	13.039	172	2164635	69.873	ng	0.00
79) Terphenyl-d14	19.815	244	4115535	70.817	ng	0.00

Target Compounds	Qvalue

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

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