

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM082522\
 Data File : BM036407.D
 Acq On : 25 Aug 2022 15:37
 Operator : CG/JU
 Sample : N4264-02RE
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 U-3RE

Quant Time: Aug 26 03:38:41 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM080122.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Aug 25 15:21:00 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.781	152	227489	20.000	ng	0.00
21) Naphthalene-d8	10.580	136	924964	20.000	ng	0.00
39) Acenaphthene-d10	14.427	164	533474	20.000	ng	0.00
64) Phenanthrene-d10	17.168	188	1016686	20.000	ng	0.00
76) Chrysene-d12	21.356	240	891977	20.000	ng	0.00
86) Perylene-d12	23.680	264	1029963	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.357	112	571995	40.765	ng	0.00
7) Phenol-d6	6.963	99	427585	23.949	ng	0.00
23) Nitrobenzene-d5	8.951	82	1553386	94.008	ng	0.00
42) 2,4,6-Tribromophenol	15.921	330	533917	85.313	ng	0.00
45) 2-Fluorobiphenyl	13.051	172	3192202	88.349	ng	0.00
79) Terphenyl-d14	19.803	244	3977662	87.959	ng	0.00

Target Compounds	Qvalue

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

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