

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM080322\
 Data File : BM036088.D
 Acq On : 03 Aug 2022 21:20
 Operator : CG/JU
 Sample : N3930-09
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 RVE-101

Quant Time: Aug 04 01:32:10 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM080122.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Aug 01 17:50:14 2022
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.863	152	178500	20.000	ng	0.00
21) Naphthalene-d8	10.669	136	870245	20.000	ng	0.00
39) Acenaphthene-d10	14.504	164	573966	20.000	ng	0.00
64) Phenanthrene-d10	17.245	188	1278626	20.000	ng	0.00
76) Chrysene-d12	21.421	240	1478254	20.000	ng	0.00
86) Perylene-d12	23.786	264	1654491	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.428	112	881476	80.062	ng	0.00
7) Phenol-d6	7.034	99	1217189	86.884	ng	0.00
23) Nitrobenzene-d5	9.028	82	769331	49.486	ng	-0.01
42) 2,4,6-Tribromophenol	15.986	330	620181	92.105	ng	-0.01
45) 2-Fluorobiphenyl	13.128	172	1954413	50.275	ng	-0.01
79) Terphenyl-d14	19.868	244	3767960	50.276	ng	0.00

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

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