

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM080522\
 Data File : BM036117.D
 Acq On : 05 Aug 2022 13:14
 Operator : CG/JU
 Sample : N3960-18
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 RVE-165

Quant Time: Aug 06 01:11:39 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM080122.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Aug 05 03:36:25 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.851	152	340276	20.000	ng	0.00
21) Naphthalene-d8	10.657	136	1443662	20.000	ng	0.00
39) Acenaphthene-d10	14.492	164	910422	20.000	ng	0.00
64) Phenanthrene-d10	17.233	188	1790547	20.000	ng	0.00
76) Chrysene-d12	21.415	240	1434979	20.000	ng	0.00
86) Perylene-d12	23.768	264	1198416	20.000	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.422	112	2334475	111.227	ng	0.00
7) Phenol-d6	7.034	99	3040481	113.849	ng	0.00
23) Nitrobenzene-d5	9.022	82	1840488	71.364	ng	0.00
42) 2,4,6-Tribromophenol	15.980	330	1192498	111.652	ng	0.00
45) 2-Fluorobiphenyl	13.121	172	4295306	69.659	ng	0.00
79) Terphenyl-d14	19.862	244	6032705	82.922	ng	0.00

Target Compounds	Qvalue

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

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