

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM080522\
 Data File : BM036130.D
 Acq On : 05 Aug 2022 21:50
 Operator : CG/JU
 Sample : N3960-16
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 RVE-163

Quant Time: Aug 06 01:13:50 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	7.857	152	273808	20.000	ng	0.00
21) Naphthalene-d8	10.657	136	1101600	20.000	ng	0.00
39) Acenaphthene-d10	14.498	164	631925	20.000	ng	0.00
64) Phenanthrene-d10	17.239	188	1232722	20.000	ng	0.00
76) Chrysene-d12	21.415	240	1122551	20.000	ng	0.00
86) Perylene-d12	23.774	264	1197040	20.000	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.446	112	852789	50.495	ng	0.02
7) Phenol-d6	7.063	99	1075935	50.068	ng	0.03
23) Nitrobenzene-d5	9.022	82	701034	35.622	ng	0.00
42) 2,4,6-Tribromophenol	15.986	330	425970	57.460	ng	0.00
45) 2-Fluorobiphenyl	13.122	172	1616009	37.758	ng	0.00
79) Terphenyl-d14	19.862	244	2284679	40.144	ng	0.00
Target Compounds						
78) Pyrene	19.657	202	195852	2.814	ng	Qvalue 100

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

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