

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM080922\
 Data File : BM036180.D
 Acq On : 09 Aug 2022 18:10
 Operator : CG/JU
 Sample : N4061-04
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 COMP1-0804

Quant Time: Aug 10 04:07:58 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 08 23:36:56 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.846	152	239738	20.000	ng	0.00
21) Naphthalene-d8	10.645	136	926102	20.000	ng	0.00
39) Acenaphthene-d10	14.486	164	489044	20.000	ng	0.00
64) Phenanthrene-d10	17.227	188	880402	20.000	ng	0.00
76) Chrysene-d12	21.409	240	873154	20.000	ng	-0.01
86) Perylene-d12	23.762	264	1052105	20.000	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.434	112	1820368	123.105	ng	0.00
7) Phenol-d6	7.051	99	2044862	108.679	ng	-0.01
23) Nitrobenzene-d5	9.016	82	1459385	88.210	ng	0.00
42) 2,4,6-Tribromophenol	15.980	330	746407	130.101	ng	0.00
45) 2-Fluorobiphenyl	13.110	172	2991444	90.315	ng	-0.01
79) Terphenyl-d14	19.857	244	3680525	83.143	ng	0.00

Target Compounds	Qvalue

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

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