

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM080822\
 Data File : BM036144.D
 Acq On : 08 Aug 2022 13:41
 Operator : CG/JU
 Sample : N4063-03
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 P-5C

Quant Time: Aug 08 23:32:04 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	425641	20.000	ng	0.00
21) Naphthalene-d8	10.657	136	1835755	20.000	ng	0.00
39) Acenaphthene-d10	14.492	164	1143627	20.000	ng	0.00
64) Phenanthrene-d10	17.239	188	2305174	20.000	ng	0.00
76) Chrysene-d12	21.427	240	1724789	20.000	ng	0.00
86) Perylene-d12	23.791	264	1413119	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.446	112	3241790	123.479	ng	0.02
7) Phenol-d6	7.069	99	4165716	124.700	ng	0.04
23) Nitrobenzene-d5	9.022	82	2529191	77.122	ng	0.00
42) 2,4,6-Tribromophenol	15.992	330	1688174	125.830	ng	0.00
45) 2-Fluorobiphenyl	13.121	172	5792500	74.784	ng	0.00
79) Terphenyl-d14	19.868	244	8196826	93.738	ng	0.00

Target Compounds	Qvalue

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

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