

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM080522\
 Data File : BM036120.D
 Acq On : 05 Aug 2022 15:07
 Operator : CG/JU
 Sample : N4064-06
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SB-14

Quant Time: Aug 06 01:12:15 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	7.852	152	310755	20.000	ng	0.00
21) Naphthalene-d8	10.657	136	1356538	20.000	ng	0.00
39) Acenaphthene-d10	14.492	164	894517	20.000	ng	0.00
64) Phenanthrene-d10	17.233	188	1882332	20.000	ng	0.00
76) Chrysene-d12	21.415	240	1719379	20.000	ng	0.00
86) Perylene-d12	23.768	264	1428938	20.000	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.422	112	1923654	100.360	ng	0.00
7) Phenol-d6	7.028	99	2536133	103.986	ng	0.00
23) Nitrobenzene-d5	9.022	82	1536425	63.400	ng	0.00
42) 2,4,6-Tribromophenol	15.981	330	1074328	102.377	ng	0.00
45) 2-Fluorobiphenyl	13.122	172	3612844	59.633	ng	0.00
79) Terphenyl-d14	19.863	244	5677233	65.128	ng	0.00

Target Compounds	Qvalue

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

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