

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM081122\
 Data File : BM036212.D
 Acq On : 11 Aug 2022 16:54
 Operator : CG/JU
 Sample : N4102-04
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 FILTER-DRUM-0808

Quant Time: Aug 12 02:16:15 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM080122.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Aug 10 22:19:26 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.828	152	227876	20.000	ng	0.00
21) Naphthalene-d8	10.628	136	950383	20.000	ng	-0.01
39) Acenaphthene-d10	14.469	164	612996	20.000	ng	0.00
64) Phenanthrene-d10	17.210	188	1416568	20.000	ng	0.00
76) Chrysene-d12	21.392	240	1522522	20.000	ng	0.00
86) Perylene-d12	23.739	264	1559632	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.410	112	1945457	138.413	ng	0.00
7) Phenol-d6	7.022	99	2266001	126.701	ng	-0.01
23) Nitrobenzene-d5	8.998	82	1623175	95.604	ng	0.00
42) 2,4,6-Tribromophenol	15.963	330	1281599	178.216	ng	0.00
45) 2-Fluorobiphenyl	13.098	172	3635912	87.575	ng	0.00
79) Terphenyl-d14	19.839	244	7433208	96.298	ng	0.00
Target Compounds						
32) Benzoic acid	9.963	122	112676	12.189	ng	Qvalue 91

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

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