

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM103122\
 Data File : BM037300.D
 Acq On : 31 Oct 2022 21:21
 Operator : CG/JU
 Sample : N5357-02
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 5051-2

Quant Time: Nov 01 05:06:01 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM102822.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Oct 31 01:34:15 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.869	152	266442	20.000	ng	0.00
21) Naphthalene-d8	10.669	136	998346	20.000	ng	0.00
39) Acenaphthene-d10	14.504	164	526247	20.000	ng	0.00
64) Phenanthrene-d10	17.245	188	1014763	20.000	ng	0.00
76) Chrysene-d12	21.415	240	1162592	20.000	ng	0.00
86) Perylene-d12	23.774	264	1329112	20.000	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.434	112	972191	63.038	ng	0.00
7) Phenol-d6	7.040	99	683207	32.640	ng	0.00
23) Nitrobenzene-d5	9.040	82	1699513	85.887	ng	0.00
42) 2,4,6-Tribromophenol	15.992	330	887888	143.170	ng	0.00
45) 2-Fluorobiphenyl	13.128	172	3574448	90.249	ng	-0.01
79) Terphenyl-d14	19.862	244	5181789	79.928	ng	0.00

Target Compounds	Qvalue

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

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