

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM103122\
 Data File : BM037298.D
 Acq On : 31 Oct 2022 20:09
 Operator : CG/JU
 Sample : N5336-06
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SB-02-16.5-17.0

Quant Time: Nov 01 05:05:28 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	249274	20.000	ng	0.00
21) Naphthalene-d8	10.669	136	924330	20.000	ng	0.00
39) Acenaphthene-d10	14.504	164	471162	20.000	ng	0.00
64) Phenanthrene-d10	17.245	188	897416	20.000	ng	0.00
76) Chrysene-d12	21.415	240	986379	20.000	ng	0.00
86) Perylene-d12	23.774	264	1169602	20.000	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.440	112	1657457	114.873	ng	0.00
7) Phenol-d6	7.040	99	2014040	102.848	ng	0.00
23) Nitrobenzene-d5	9.039	82	1184276	64.641	ng	0.00
42) 2,4,6-Tribromophenol	15.992	330	594973	107.154	ng	0.00
45) 2-Fluorobiphenyl	13.127	172	2299530	64.848	ng	-0.01
79) Terphenyl-d14	19.862	244	3112653	56.589	ng	0.00

Target Compounds Qvalue

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

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