

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM103122\
 Data File : BM037297.D
 Acq On : 31 Oct 2022 19:33
 Operator : CG/JU
 Sample : N5337-04
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SB-02-COMP

Quant Time: Nov 01 05:05:08 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	275914	20.000	ng	0.00
21) Naphthalene-d8	10.669	136	1081509	20.000	ng	0.00
39) Acenaphthene-d10	14.504	164	601244	20.000	ng	0.00
64) Phenanthrene-d10	17.245	188	1149038	20.000	ng	0.00
76) Chrysene-d12	21.415	240	1207958	20.000	ng	0.00
86) Perylene-d12	23.780	264	1346147	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.440	112	1777550	111.302	ng	0.00
7) Phenol-d6	7.040	99	2212519	102.075	ng	0.00
23) Nitrobenzene-d5	9.040	82	1315838	61.384	ng	0.00
42) 2,4,6-Tribromophenol	15.992	330	755288	106.597	ng	0.00
45) 2-Fluorobiphenyl	13.128	172	2812311	62.149	ng	-0.01
79) Terphenyl-d14	19.862	244	4109161	61.002	ng	0.00
Target Compounds						
71) Phenanthrene	17.286	178	199238	2.795	ng	Qvalue # 92

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

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