

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM112122\
 Data File : BM037647.D
 Acq On : 22 Nov 2022 18:15
 Operator : CG/JU
 Sample : N5577-13
 Misc :
 ALS Vial : 47 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C9837

Quant Time: Nov 22 22:52:34 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-SIM-BM111622.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Nov 21 22:36:32 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.129	152	8675	0.400	ng/ul	0.00
4) Naphthalene-d8	10.946	136	27014	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.746	164	16300	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.482	188	35313	0.400	ng/ul	0.00
17) Chrysene-d12	21.637	240	27304	0.400	ng/ul	# 0.00
23) Perylene-d12	24.124	264	27638	0.400	ng/ul	# 0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.450	96	51807	5.136	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.524	152	12273	0.274	ng/ul	0.00
18) Fluoranthene-d10	19.493	212	30595	0.339	ng/ul	0.00
Target Compounds						Qvalue
5) Naphthalene	10.996	128	2524	0.031	ng/ul#	89
7) 2-Methylnaphthalene	12.596	142	1686	0.030	ng/ul	100
8) 1-Methylnaphthalene	12.810	142	1206	0.021	ng/ul	100
15) Phenanthrene	17.520	178	3154	0.026	ng/ul#	86
19) Fluoranthene	19.521	202	2699	0.020	ng/ul#	67

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

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