

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM102122\
 Data File : BM037161.D
 Acq On : 21 Oct 2022 06:20
 Operator : CG/JU
 Sample : N5103-16
 Misc :
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C0BA1

Quant Time: Oct 21 06:50:52 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-SIM-BM101922.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Oct 20 02:11:25 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.905	152	4588	0.400	ng/ul	0.00
4) Naphthalene-d8	10.704	136	14391	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.533	164	8473	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.275	188	17716	0.400	ng/ul	0.00
17) Chrysene-d12	21.450	240	14671	0.400	ng/ul	# 0.00
23) Perylene-d12	23.823	264	14542	0.400	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.314	96	20728	3.705	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.294	152	5995	0.293	ng/ul	0.00
18) Fluoranthene-d10	19.298	212	17084	0.385	ng/ul	0.00
Target Compounds						
					Qvalue	
5) Naphthalene	10.754	128	1031	0.024	ng/ul#	81
7) 2-Methylnaphthalene	12.365	142	579	0.023	ng/ul	91
15) Phenanthrene	17.313	178	2054	0.034	ng/ul#	88
19) Fluoranthene	19.326	202	3952	0.062	ng/ul#	90
20) Pyrene	19.689	202	3768	0.058	ng/ul#	84
21) Benzo(a)anthracene	21.435	228	2314	0.041	ng/ul#	89
22) Chrysene	21.488	228	2260	0.034	ng/ul#	89
24) Benzo(b)fluoranthene	23.101	252	3289	0.043	ng/ul#	1
26) Benzo(a)pyrene	23.718	252	2278	0.037	ng/ul#	1

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

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