

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM101022\
 Data File : BM037004.D
 Acq On : 13 Oct 2022 03:54
 Operator : CG/JU
 Sample : N4976-08
 Misc :
 ALS Vial : 109 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 CB8Y5

Quant Time: Oct 13 05:51:10 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-SIM-BM100622.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Oct 11 09:03:52 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.947	152	8421	0.400	ng/ul	0.00
4) Naphthalene-d8	10.748	136	30411	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.570	164	19042	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.309	188	34917	0.400	ng/ul	# 0.00
17) Chrysene-d12	21.479	240	20730	0.400	ng/ul	# 0.00
23) Perylene-d12	23.876	264	17043	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.344	96	37282	3.139	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.332	152	10899	0.237	ng/ul	0.00
18) Fluoranthene-d10	19.331	212	24798	0.400	ng/ul	0.00
Target Compounds						
					Qvalue	
5) Naphthalene	10.798	128	1855	0.021	ng/ul#	48
10) Acenaphthylene	14.311	152	3094	0.039	ng/ul#	6
11) Acenaphthene	14.635	153	41360	0.602	ng/ul	99
12) Fluorene	15.616	166	2904	0.037	ng/ul#	67
15) Phenanthrene	17.351	178	2474	0.022	ng/ul#	8
16) Anthracene	17.444	178	6224	0.067	ng/ul#	53

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

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM101022\
Data File : BM037004.D
Acq On : 13 Oct 2022 03:54
Operator : CG/JU
Sample : N4976-08
Misc :
ALS Vial : 109 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
CB8Y5

Quant Time: Oct 13 05:51:10 2022
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-SIM-BM100622.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Oct 11 09:03:52 2022
Response via : Initial Calibration

