

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM080921\
 Data File : BM031603.D
 Acq On : 09 Aug 2021 14:55
 Operator : CG/JU
 Sample : M3244-04
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 JCE04

Quant Time: Aug 10 02:47:47 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-SIM-BM072821.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Aug 09 10:16:43 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.580	152	1119	0.400	ng/ul	0.00
4) Naphthalene-d8	10.338	136	4122	0.400	ng/ul	# 0.00
9) Acenaphthene-d10	14.212	164	2501	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.961	188	4865	0.400	ng/ul	# 0.00
17) Chrysene-d12	21.159	240	3801	0.400	ng/ul	# 0.00
23) Perylene-d12	23.314	264	3563	0.400	ng/ul	# 0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.186	96	3640	2.929	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.935	152	1723	0.248	ng/ul	0.00
18) Fluoranthene-d10	18.993	212	3442	0.271	ng/ul	0.00
Target Compounds						
					Qvalue	
5) Naphthalene	10.388	128	357	0.029	ng/ul#	56
15) Phenanthrene	17.002	178	514	0.033	ng/ul#	64
19) Fluoranthene	19.027	202	785	0.048	ng/ul#	1
20) Pyrene	19.390	202	562	0.034	ng/ul#	1
22) Chrysene	21.197	228	1219	0.077	ng/ul#	79
24) Benzo(b)fluoranthene	22.679	252	806	0.050	ng/ul#	1
25) Benzo(k)fluoranthene	22.730	252	403	0.024	ng/ul#	1

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

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM080921\
Data File : BM031603.D
Acq On : 09 Aug 2021 14:55
Operator : CG/JU
Sample : M3244-04
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
JCE04

Quant Time: Aug 10 02:47:47 2021
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-SIM-BM072821.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Aug 09 10:16:43 2021
Response via : Initial Calibration

