

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM080921\
 Data File : BM031613.D
 Acq On : 09 Aug 2021 21:07
 Operator : CG/JU
 Sample : M3244-03
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 JCE03

Quant Time: Aug 10 02:55:15 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 : Tue Aug 10 02:43:39 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.580	152	1163	0.400	ng/ul	0.00
4) Naphthalene-d8	10.343	136	4249	0.400	ng/ul #	0.00
9) Acenaphthene-d10	14.212	164	2375	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.961	188	4582	0.400	ng/ul #	0.00
17) Chrysene-d12	21.156	240	3711	0.400	ng/ul #	0.00
23) Perylene-d12	23.311	264	3766	0.400	ng/ul #	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.199	96	34	0.026	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.935	152	917	0.128	ng/ul	0.00
18) Fluoranthene-d10	18.997	212	1694	0.137	ng/ul	0.00
Target Compounds						
					Qvalue	
5) Naphthalene	10.392	128	357	0.028	ng/ul#	60
15) Phenanthrene	17.002	178	545	0.037	ng/ul#	75
19) Fluoranthene	19.027	202	640	0.040	ng/ul#	56
20) Pyrene	19.390	202	544	0.034	ng/ul#	41
22) Chrysene	21.193	228	574	0.037	ng/ul#	61
24) Benzo(b)fluoranthene	22.672	252	697	0.041	ng/ul#	1

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

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