

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM080221\
 Data File : BM031437.D
 Acq On : 04 Aug 2021 07:21
 Operator : CG/JU
 Sample : M3122-21
 Misc :
 ALS Vial : 67 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 BGET4

Manual Integrations
 APPROVED

mohammad
 8/4/2021 3:48:56 PM

Quant Time: Aug 04 08:21:18 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 : Wed Aug 04 04:12:19 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.594	152	1119	0.400	ng/ul	0.00
4) Naphthalene-d8	10.356	136	4310	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.224	164	2596	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.971	188	4896	0.400	ng/ul	# 0.00
17) Chrysene-d12	21.166	240	3627	0.400	ng/ul	0.00
23) Perylene-d12	23.324	264	3595	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.192	96	3382	2.721	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.953	152	1203	0.166	ng/ul	0.00
18) Fluoranthene-d10	19.005	212	2381	0.197	ng/ul	0.00
Target Compounds						
					Qvalue	
15) Phenanthrene	17.013	178	475	0.031	ng/ul#	76
19) Fluoranthene	19.035	202	1699	0.110	ng/ul#	88
20) Pyrene	19.397	202	770	0.049	ng/ul#	73
21) Benzo(a)anthracene	21.149	228	689	0.047	ng/ul#	80
22) Chrysene	21.202	228	942	0.062	ng/ul#	88
24) Benzo(b)fluoranthene	22.682	252	1397m	0.086	ng/ul	
25) Benzo(k)fluoranthene	22.725	252	578m	0.035	ng/ul	

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

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