

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM080221\
 Data File : BM031433.D
 Acq On : 04 Aug 2021 04:55
 Operator : CG/JU
 Sample : M3122-22
 Misc :
 ALS Vial : 63 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 BGET9

Manual Integrations
 APPROVED

mohammad
 8/4/2021 3:48:42 PM

Quant Time: Aug 04 05:27:50 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	1062	0.400	ng/ul	0.00
4) Naphthalene-d8	10.356	136	4022	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.224	164	2407	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.975	188	4496	0.400	ng/ul	0.00
17) Chrysene-d12	21.166	240	3285	0.400	ng/ul	0.00
23) Perylene-d12	23.324	264	3257	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.193	96	2959	2.509	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.953	152	1042	0.154	ng/ul	0.00
18) Fluoranthene-d10	19.005	212	1884	0.172	ng/ul	0.00
Target Compounds						
					Qvalue	
15) Phenanthrene	17.016	178	354	0.025	ng/ul#	67
19) Fluoranthene	19.035	202	1026	0.073	ng/ul#	80
20) Pyrene	19.397	202	435	0.031	ng/ul#	53
21) Benzo(a)anthracene	21.152	228	391	0.029	ng/ul#	77
22) Chrysene	21.202	228	574	0.042	ng/ul#	80
24) Benzo(b)fluoranthene	22.684	252	766m	0.052	ng/ul	

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

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