

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM080221\
 Data File : BM031439.D
 Acq On : 04 Aug 2021 08:34
 Operator : CG/JU
 Sample : M3122-20
 Misc :
 ALS Vial : 69 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 BGEW9

Manual Integrations
 APPROVED

mohammad
 8/4/2021 3:48:59 PM

Quant Time: Aug 04 09:21:51 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 03 13:47:07 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.594	152	1031	0.400	ng/ul	0.00
4) Naphthalene-d8	10.356	136	3774	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.224	164	2232	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.975	188	4396	0.400	ng/ul	0.00
17) Chrysene-d12	21.169	240	3491	0.400	ng/ul	0.00
23) Perylene-d12	23.329	264	3424	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.193	96	3018	2.636	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.953	152	1002	0.158	ng/ul	0.00
18) Fluoranthene-d10	19.005	212	1883	0.162	ng/ul	0.00
Target Compounds						
15) Phenanthrene	17.017	178	386	0.028	ng/ul#	71
19) Fluoranthene	19.035	202	1106	0.074	ng/ul#	82
20) Pyrene	19.398	202	528m	0.035	ng/ul	
21) Benzo(a)anthracene	21.154	228	440	0.031	ng/ul#	72
22) Chrysene	21.205	228	606	0.042	ng/ul#	83
24) Benzo(b)fluoranthene	22.687	252	879m	0.057	ng/ul	
25) Benzo(k)fluoranthene	22.723	252	405m	0.025	ng/ul	

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

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