

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM080221\
 Data File : BM031434.D
 Acq On : 04 Aug 2021 05:32
 Operator : CG/JU
 Sample : M3122-12
 Misc :
 ALS Vial : 64 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 BGEQ8

Manual Integrations
 APPROVED

mohammad
 8/4/2021 3:48:44 PM

Quant Time: Aug 04 06:04:53 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	7.597	152	1173	0.400	ng/ul	0.00
4) Naphthalene-d8	10.355	136	4372	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.227	164	2635	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.974	188	4998	0.400	ng/ul	# 0.00
17) Chrysene-d12	21.166	240	3858	0.400	ng/ul	0.00
23) Perylene-d12	23.323	264	3830	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.192	96	2995	2.299	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.952	152	1136	0.154	ng/ul	0.00
18) Fluoranthene-d10	19.008	212	2300	0.179	ng/ul	0.00
Target Compounds						
					Qvalue	
10) Acenaphthylene	13.944	152	257	0.023	ng/ul#	55
12) Fluorene	15.281	166	258	0.024	ng/ul#	87
15) Phenanthrene	17.016	178	3569	0.225	ng/ul	99
16) Anthracene	17.106	178	1370	0.092	ng/ul#	90
19) Fluoranthene	19.035	202	7972	0.484	ng/ul	98
20) Pyrene	19.397	202	4266m	0.255	ng/ul	
21) Benzo(a)anthracene	21.151	228	4727	0.303	ng/ul	95
22) Chrysene	21.202	228	4372	0.271	ng/ul	97
24) Benzo(b)fluoranthene	22.684	252	5163m	0.299	ng/ul	
25) Benzo(k)fluoranthene	22.720	252	1873m	0.105	ng/ul	
26) Benzo(a)pyrene	23.227	252	1408	0.093	ng/ul#	48
27) Indeno(1,2,3-cd)pyrene	25.458	276	1104	0.057	ng/ul#	94
28) Dibenzo(a,h)anthracene	25.472	278	662	0.042	ng/ul#	34

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

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