

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM080921\
 Data File : BM031630.D
 Acq On : 10 Aug 2021 08:07
 Operator : CG/JU
 Sample : M3169-08
 Misc :
 ALS Vial : 38 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C00A7

Manual Integrations
 APPROVED

mohammad
 8/10/2021 4:37:13 PM

Quant Time: Aug 10 09:20:02 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 08:05:02 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.576	152	1403	0.400	ng/ul	0.00
4) Naphthalene-d8	10.338	136	5067	0.400	ng/ul	# 0.00
9) Acenaphthene-d10	14.208	164	3609	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.961	188	5423	0.400	ng/ul	# 0.00
17) Chrysene-d12	21.156	240	4336	0.400	ng/ul	# 0.00
23) Perylene-d12	23.314	264	4805	0.400	ng/ul	# 0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.182	96	7878	5.056	ng/ul	-0.01
6) 2-Methylnaphthalene-d10	11.930	152	3164	0.371	ng/ul	0.00
18) Fluoranthene-d10	18.997	212	5911	0.408	ng/ul	0.00
Target Compounds						
					Qvalue	
5) Naphthalene	10.383	128	39289	2.603	ng/ul	98
7) 2-Methylnaphthalene	12.002	142	41083	3.984	ng/ul	100
8) 1-Methylnaphthalene	12.227	142	30698	3.013	ng/ul	99
10) Acenaphthylene	13.932	152	24894	1.620	ng/ul	99
11) Acenaphthene	14.273	153	4915m	0.380	ng/ul	
12) Fluorene	15.266	166	7869	0.529	ng/ul#	78
15) Phenanthrene	17.002	178	207912	12.074	ng/ul	98
16) Anthracene	17.093	178	31234	1.937	ng/ul	99
19) Fluoranthene	19.024	202	271189	14.647	ng/ul#	96
20) Pyrene	19.390	202	235953	12.565	ng/ul	96
21) Benzo(a)anthracene	21.139	228	124311	7.100	ng/ul#	88
22) Chrysene	21.193	228	191593	10.576	ng/ul#	94
24) Benzo(b)fluoranthene	22.677	252	199198m	9.201	ng/ul	
25) Benzo(k)fluoranthene	22.713	252	52849m	2.363	ng/ul	
26) Benzo(a)pyrene	23.222	252	120111m	6.335	ng/ul	
27) Indeno(1,2,3-cd)pyrene	25.453	276	82314	3.359	ng/ul#	100
28) Dibenzo(a,h)anthracene	25.458	278	26196	1.326	ng/ul#	72
29) Benzo(g,h,i)perylene	26.105	276	78297	3.745	ng/ul	98

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

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