

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM111720\
 Data File : BM027815.D
 Acq On : 17 Nov 2020 12:03
 Operator : JU/CG
 Sample : L4762-10
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C0AC6

Manual Integrations
 APPROVED

mohammad
 11/19/2020 10:32:50 AM

Quant Time: Nov 17 12:49:09 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\SOM-EPA-SIM-BM111020.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Nov 17 09:19:02 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.38	152	1181	0.40	ng/ul	0.00
2) Naphthalene-d8	11.22	136	4872	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.99	164	3192	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.70	188	6175m	0.40	ng/ul	0.00
16) Chrysene-d12	21.80	240	3385	0.40	ng/ul	0.00
20) Perylene-d12	24.25	264	2628	0.40	ng/ul	0.02
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.78	152	2383	0.26	ng/ul	0.00
14) Fluoranthene-d10	19.69	212	3482	0.19	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	11.26	128	31769	2.041	ng/ul	95
5) 2-Methylnaphthalene	12.85	142	69376	6.406	ng/ul	99
7) Acenaphthylene	14.71	152	4530	0.288	ng/ul#	65
8) Acenaphthene	15.05	153	3547	0.295	ng/ul#	63
9) Fluorene	16.02	166	5742	0.397	ng/ul#	83
12) Phenanthrene	17.74	178	83736	4.166	ng/ul	98
13) Anthracene	17.83	178	11078m	0.566	ng/ul	
15) Fluoranthene	19.72	202	34223	1.487	ng/ul	95
17) Pyrene	20.08	202	49325	3.286	ng/ul#	93
18) Benzo(a)anthracene	21.79	228	18769	1.387	ng/ul#	47
19) Chrysene	21.84	228	30407	2.207	ng/ul#	61
21) Benzo(b)fluoranthene	23.50	252	23020m	1.914	ng/ul	
22) Benzo(k)fluoranthene	23.55	252	7985m	0.676	ng/ul	
23) Benzo(a)pyrene	24.14	252	14880	1.407	ng/ul#	34
24) Indeno(1,2,3-cd)pyrene	26.77	276	11735	0.974	ng/ul#	96
25) Dibenzo(a,h)anthracene	26.77	278	3766	0.374	ng/ul#	1
26) Benzo(a,h,i)perylene	27.57	276	20699	2.030	ng/ul#	76

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

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