

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM040420\
 Data File : BM025928.D
 Acq On : 04 Apr 2020 00:48
 Operator : CG/JU
 Sample : L2062-07DL 5X
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 F3E06DL

Manual Integrations
 APPROVED

mohammad
 4/6/2020 1:41:29 PM

Quant Time: Apr 04 02:02:47 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\SOM-EPA-SIM-BM040220.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Apr 03 22:53:21 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.56	152	4447	0.40	ng/ul	0.00
2) Naphthalene-d8	10.33	136	19810	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.20	164	13096	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.95	188	30475	0.40	ng/ul	0.00
16) Chrysene-d12	21.14	240	30483	0.40	ng/ul	0.00
20) Perylene-d12	23.28	264	26945	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.93	152	1604	0.05	ng/ul	0.00
14) Fluoranthene-d10	18.98	212	4883	0.05	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.38	128	1184	0.020	ng/ul#	83
7) Acenaphthylene	13.92	152	19598	0.350	ng/ul	99
8) Acenaphthene	14.26	153	2153	0.046	ng/ul	97
9) Fluorene	15.25	166	5179	0.090	ng/ul	98
12) Phenanthrene	16.99	178	19853	0.207	ng/ul	99
13) Anthracene	17.08	178	436002	5.140	ng/ul	99
15) Fluoranthene	19.01	202	790037	6.660	ng/ul	99
17) Pyrene	19.37	202	1310147	12.001	ng/ul	99
18) Benzo(a)anthracene	21.12	228	851363	7.813	ng/ul	99
19) Chrysene	21.18	228	1309946	11.014	ng/ul	99
21) Benzo(b)fluoranthene	22.65	252	1260636m	12.713	ng/ul	
22) Benzo(k)fluoranthene	22.69	252	418317m	3.934	ng/ul	
23) Benzo(a)pyrene	23.19	252	530097	5.982	ng/ul	96
24) Indeno(1,2,3-cd)pyrene	25.40	276	192247	1.620	ng/ul#	100
25) Dibenzo(a,h)anthracene	25.40	278	58021	0.586	ng/ul	98
26) Benzo(a,h,i)perylene	26.04	276	101439	1.010	ng/ul	99

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

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