

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM030320\
 Data File : BM025215.D
 Acq On : 03 Mar 2020 16:57
 Operator : CG/JU
 Sample : L1507-07
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 DBD05

Manual Integrations
 APPROVED

mohammad
 3/4/2020 7:52:50 AM

Quant Time: Mar 03 17:31:23 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\SOM-EPA-SIM-BM030320.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Mar 03 14:30:08 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.68	152	2828	0.40	ng/ul	0.00
2) Naphthalene-d8	10.45	136	11154	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.31	164	7899	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.05	188	19175	0.40	ng/ul	0.00
16) Chrysene-d12	21.24	240	19473	0.40	ng/ul	0.00
20) Perylene-d12	23.43	264	18972	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.04	152	3486	0.17	ng/ul	0.00
14) Fluoranthene-d10	19.08	212	10687	0.17	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.50	128	2330	0.069	ng/ul#	90
5) 2-Methylnaphthalene	12.12	142	870	0.036	ng/ul	97
7) Acenaphthylene	14.02	152	1105	0.030	ng/ul	95
12) Phenanthrene	17.09	178	3382	0.055	ng/ul	98
13) Anthracene	17.18	178	1477m	0.027	ng/ul	
15) Fluoranthene	19.11	202	30481	0.394	ng/ul	95
17) Pyrene	19.47	202	28637	0.388	ng/ul	95
18) Benzo(a)anthracene	21.22	228	26754	0.374	ng/ul	99
19) Chrysene	21.27	228	13716m	0.174	ng/ul	
21) Benzo(b)fluoranthene	22.78	252	23386m	0.308	ng/ul	
22) Benzo(k)fluoranthene	22.82	252	10482m	0.132	ng/ul	
23) Benzo(a)pyrene	23.33	252	8766	0.138	ng/ul#	92
24) Indeno(1,2,3-cd)pyrene	25.61	276	4334	0.055	ng/ul#	100
25) Dibenzo(a,h)anthracene	25.62	278	1533	0.023	ng/ul#	59
26) Benzo(g,h,i)perylene	26.28	276	2657	0.039	ng/ul	95

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

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