

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM022820\
 Data File : BM025138.D
 Acq On : 28 Feb 2020 18:08
 Operator : CG/JU
 Sample : L1507-04
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 DBD02

Manual Integrations
 APPROVED

Sohil
 3/3/2020 8:38:36 AM

Quant Time: Feb 29 05:23:59 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\SOM-EPA-SIM-BM022420.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Feb 29 04:43:59 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.69	152	2848	0.40	ng/ul	0.00
2) Naphthalene-d8	10.46	136	10634	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.31	164	6484	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.06	188	14894	0.40	ng/ul	0.00
16) Chrysene-d12	21.24	240	14932	0.40	ng/ul	0.00
20) Perylene-d12	23.44	264	16331	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.05	152	396	0.02	ng/ul	0.00
14) Fluoranthene-d10	19.08	212	1075	0.02	ng/ul	0.00
Target Compounds				Ovalue		
12) Phenanthrene	17.10	178	1095	0.021	ng/ul#	87
13) Anthracene	17.19	178	1100	0.022	ng/ul#	82
15) Fluoranthene	19.11	202	5157	0.068	ng/ul	99
17) Pyrene	19.48	202	8508	0.126	ng/ul	97
18) Benzo(a)anthracene	21.22	228	6004	0.093	ng/ul	96
19) Chrysene	21.26	228	9219	0.144	ng/ul	99
21) Benzo(b)fluoranthene	22.79	252	15608m	0.244	ng/ul	
22) Benzo(k)fluoranthene	22.83	252	6249m	0.099	ng/ul	
23) Benzo(a)pyrene	23.34	252	7193	0.123	ng/ul	93
24) Indeno(1,2,3-cd)pyrene	25.63	276	4051	0.055	ng/ul#	100
26) Benzo(g,h,i)perylene	26.29	276	2419	0.039	ng/ul#	92

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

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