

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM022820\
 Data File : BM025139.D
 Acq On : 28 Feb 2020 18:44
 Operator : CG/JU
 Sample : L1507-05
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 DBD03

Manual Integrations
 APPROVED

Sohil
 3/3/2020 8:38:37 AM

Quant Time: Feb 29 05:26:08 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	2375	0.40	ng/ul	0.00
2) Naphthalene-d8	10.46	136	9093	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.31	164	6199	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.06	188	13999m	0.40	ng/ul	0.00
16) Chrysene-d12	21.24	240	14611	0.40	ng/ul	0.00
20) Perylene-d12	23.44	264	16460	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.05	152	7574	0.44	ng/ul	0.00
14) Fluoranthene-d10	19.08	212	20710	0.43	ng/ul	0.00
Target Compounds					Ovalue	
3) Naphthalene	10.51	128	1865	0.067	ng/ul#	92
8) Acenaphthene	14.38	153	6948	0.289	ng/ul	99
9) Fluorene	15.37	166	7842	0.260	ng/ul	95
12) Phenanthrene	17.10	178	4154	0.084	ng/ul	99
13) Anthracene	17.19	178	1425	0.030	ng/ul#	89
15) Fluoranthene	19.11	202	5062	0.071	ng/ul	99
17) Pyrene	19.48	202	3908	0.059	ng/ul#	91
18) Benzo(a)anthracene	21.22	228	1694	0.027	ng/ul#	90
19) Chrysene	21.27	228	2563	0.041	ng/ul#	90
21) Benzo(b)fluoranthene	22.79	252	4288m	0.066	ng/ul	
22) Benzo(k)fluoranthene	22.82	252	1519m	0.024	ng/ul	
23) Benzo(a)pyrene	23.34	252	1303	0.022	ng/ul#	61
24) Indeno(1,2,3-cd)pyrene	25.63	276	1610	0.022	ng/ul#	100

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

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