

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM030920\
 Data File : BM025464.D
 Acq On : 10 Mar 2020 22:04
 Operator : CG/JU
 Sample : L1617-14
 Misc :
 ALS Vial : 58 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :

Quant Time: Mar 11 00:12:19 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\SOM-EPA-SIM-BM030920.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Mar 10 23:48:01 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.66	152	2676	0.40	ng/ul	0.00
2) Naphthalene-d8	10.43	136	11644	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.29	164	8435	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.03	188	19081	0.40	ng/ul	0.00
16) Chrysene-d12	21.22	240	13994	0.40	ng/ul	0.00
20) Perylene-d12	23.40	264	12696	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.02	152	7718	0.37	ng/ul	0.00
14) Fluoranthene-d10	19.06	212	21080	0.40	ng/ul	0.00
Target Compounds				Ovalue		
12) Phenanthrene	17.07	178	27315	0.452	ng/ul	99
13) Anthracene	17.16	178	7834	0.149	ng/ul	97
15) Fluoranthene	19.09	202	87708	1.338	ng/ul	98
17) Pyrene	19.45	202	59494	0.947	ng/ul	100
18) Benzo(a)anthracene	21.20	228	12942	0.259	ng/ul	99
19) Chrysene	21.26	228	12642	0.225	ng/ul	98
21) Benzo(b)fluoranthene	22.76	252	9305m	0.185	ng/ul	
22) Benzo(k)fluoranthene	22.79	252	2880m	0.055	ng/ul	
23) Benzo(a)pyrene	23.31	252	2938	0.070	ng/ul#	66
24) Indeno(1,2,3-cd)pyrene	25.58	276	1600	0.028	ng/ul#	63

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

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