

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM030920\
 Data File : BM025468.D
 Acq On : 11 Mar 2020 00:31
 Operator : CG/JU
 Sample : L1617-09
 Misc :
 ALS Vial : 62 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :

Quant Time: Mar 11 02:31:42 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	2414	0.40	ng/ul	0.00
2) Naphthalene-d8	10.43	136	9811	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.29	164	7093	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.03	188	15721	0.40	ng/ul	0.00
16) Chrysene-d12	21.22	240	10904	0.40	ng/ul	0.00
20) Perylene-d12	23.40	264	11086	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.02	152	5686	0.32	ng/ul	0.00
14) Fluoranthene-d10	19.06	212	14224	0.33	ng/ul	0.00
Target Compounds				Ovalue		
12) Phenanthrene	17.07	178	1345	0.027	ng/ul#	87
13) Anthracene	17.16	178	957	0.022	ng/ul#	81
15) Fluoranthene	19.09	202	1771	0.033	ng/ul	87
17) Pyrene	19.45	202	2316m	0.047	ng/ul	
18) Benzo(a)anthracene	21.20	228	796	0.020	ng/ul#	78
19) Chrysene	21.25	228	1160	0.026	ng/ul#	89
21) Benzo(b)fluoranthene	22.76	252	1544m	0.035	ng/ul	
22) Benzo(k)fluoranthene	22.79	252	1074m	0.023	ng/ul	
23) Benzo(a)pyrene	23.30	252	787	0.022	ng/ul#	1
24) Indeno(1,2,3-cd)pyrene	25.58	276	1170	0.024	ng/ul#	63
26) Benzo(g,h,i)perylene	26.24	276	1249	0.029	ng/ul#	72

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

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