

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM030620\
 Data File : BM025387.D
 Acq On : 08 Mar 2020 16:11
 Operator : CG/JU
 Sample : L1617-12
 Misc :
 ALS Vial : 73 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 DBDS6

Manual Integrations
 APPROVED

mohammad
 3/9/2020 3:03:31 PM

Quant Time: Mar 09 02:08:39 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 : Wed Mar 04 08:02:03 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.67	152	4337	0.40	ng/ul	-0.01
2) Naphthalene-d8	10.43	136	18558	0.40	ng/ul	-0.01
6) Acenaphthene-d10	14.29	164	13185	0.40	ng/ul	-0.01
10) Phenanthrene-d10	17.04	188	30951	0.40	ng/ul	-0.01
16) Chrysene-d12	21.22	240	25915	0.40	ng/ul	0.00
20) Perylene-d12	23.41	264	22460	0.40	ng/ul	-0.01
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.03	152	4630	0.13	ng/ul	-0.01
14) Fluoranthene-d10	19.06	212	12869	0.13	ng/ul	-0.01
Target Compounds					Ovalue	
8) Acenaphthene	14.36	153	4031	0.084	ng/ul	99
9) Fluorene	15.34	166	13640	0.227	ng/ul	96
12) Phenanthrene	17.07	178	121561	1.218	ng/ul	95
13) Anthracene	17.17	178	7640	0.086	ng/ul	98
15) Fluoranthene	19.09	202	74609	0.597	ng/ul	95
17) Pyrene	19.46	202	50481	0.514	ng/ul	96
18) Benzo(a)anthracene	21.20	228	7046	0.074	ng/ul	99
19) Chrysene	21.26	228	6444	0.062	ng/ul	99
21) Benzo(b)fluoranthene	22.76	252	3599m	0.040	ng/ul	
23) Benzo(a)pyrene	23.31	252	1705	0.023	ng/ul#	35

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

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