

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM030920\
 Data File : BM025465.D
 Acq On : 10 Mar 2020 22:41
 Operator : CG/JU
 Sample : L1617-22
 Misc :
 ALS Vial : 59 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 DBDT6

Manual Integrations
 APPROVED

mohammad
 3/11/2020 8:01:58 AM

Quant Time: Mar 11 00:14:49 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	3024	0.40	ng/ul	0.00
2) Naphthalene-d8	10.43	136	12680	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.29	164	9019	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.03	188	20195	0.40	ng/ul	0.00
16) Chrysene-d12	21.22	240	13745	0.40	ng/ul	0.00
20) Perylene-d12	23.40	264	12920	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.03	152	5059	0.22	ng/ul	0.00
14) Fluoranthene-d10	19.06	212	12270	0.22	ng/ul	0.00
Target Compounds				Ovalue		
15) Fluoranthene	19.09	202	2999	0.043	ng/ul	92
17) Pyrene	19.45	202	3408	0.055	ng/ul#	94
18) Benzo(a)anthracene	21.20	228	2485	0.051	ng/ul	93
19) Chrysene	21.24	228	2227m	0.040	ng/ul	
21) Benzo(b)fluoranthene	22.76	252	3750m	0.073	ng/ul	
22) Benzo(k)fluoranthene	22.80	252	1209m	0.023	ng/ul	
23) Benzo(a)pyrene	23.31	252	1815	0.043	ng/ul#	44

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

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