

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM030920\
 Data File : BM025457.D
 Acq On : 10 Mar 2020 17:48
 Operator : CG/JU
 Sample : L1617-19DL 5X
 Misc :
 ALS Vial : 51 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 DBDT3DL

Manual Integrations
 APPROVED

mohammad
 3/11/2020 8:01:36 AM

Quant Time: Mar 10 18:21:22 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 : Mon Mar 09 13:32:56 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.66	152	3122	0.40	ng/ul	0.00
2) Naphthalene-d8	10.43	136	13300	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.29	164	9260	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.03	188	19248	0.40	ng/ul	0.00
16) Chrysene-d12	21.22	240	13225	0.40	ng/ul	0.00
20) Perylene-d12	23.40	264	13476	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.03	152	1551	0.06	ng/ul	0.00
14) Fluoranthene-d10	19.06	212	3176	0.06	ng/ul	0.00
Target Compounds					Ovalue	
8) Acenaphthene	14.35	153	16614	0.501	ng/ul	100
9) Fluorene	15.34	166	30026	0.743	ng/ul	100
12) Phenanthrene	17.07	178	115168	1.888	ng/ul	99
13) Anthracene	17.16	178	7222	0.136	ng/ul	98
15) Fluoranthene	19.09	202	75003	1.134	ng/ul	99
17) Pyrene	19.45	202	47167	0.795	ng/ul	98
18) Benzo(a)anthracene	21.21	228	3368	0.071	ng/ul	96
19) Chrysene	21.26	228	3739	0.070	ng/ul	96
21) Benzo(b)fluoranthene	22.76	252	3174m	0.059	ng/ul	
22) Benzo(k)fluoranthene	22.80	252	1568m	0.028	ng/ul	
23) Benzo(a)pyrene	23.31	252	1535	0.035	ng/ul#	34

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

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