

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM030920\
 Data File : BM025454.D
 Acq On : 10 Mar 2020 15:58
 Operator : CG/JU
 Sample : L1617-13DL 5X
 Misc :
 ALS Vial : 48 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 DBDS7DL

Manual Integrations
 APPROVED

mohammad
 3/11/2020 8:01:30 AM

Quant Time: Mar 10 16:36:17 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	2755	0.40	ng/ul	0.00
2) Naphthalene-d8	10.43	136	11297	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.29	164	7724	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.03	188	16522	0.40	ng/ul	0.00
16) Chrysene-d12	21.22	240	10915	0.40	ng/ul	0.00
20) Perylene-d12	23.41	264	10382	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.03	152	1365	0.07	ng/ul	0.00
14) Fluoranthene-d10	19.06	212	2796	0.06	ng/ul	0.00
Target Compounds				Ovalue		
12) Phenanthrene	17.07	178	5628	0.107	ng/ul	98
15) Fluoranthene	19.09	202	44444	0.783	ng/ul	100
17) Pyrene	19.45	202	30748	0.628	ng/ul	99
18) Benzo(a)anthracene	21.20	228	7459	0.191	ng/ul	98
19) Chrysene	21.26	228	9238	0.211	ng/ul	99
21) Benzo(b)fluoranthene	22.76	252	4431m	0.108	ng/ul	
22) Benzo(k)fluoranthene	22.80	252	1967m	0.046	ng/ul	
23) Benzo(a)pyrene	23.31	252	1429	0.042	ng/ul#	40

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

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