

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM031720\
 Data File : BM025621.D
 Acq On : 17 Mar 2020 21:33
 Operator : CG/JU
 Sample : L1516-06DL 10X
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 DBD24DL

Manual Integrations
 APPROVED

mohammad
 3/18/2020 9:02:12 AM

Quant Time: Mar 18 01:53:12 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 17 17:05:05 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.59	152	2343	0.40	ng/ul	0.00
2) Naphthalene-d8	10.35	136	9007	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.22	164	5813	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.96	188	13964	0.40	ng/ul	0.00
16) Chrysene-d12	21.16	240	14263	0.40	ng/ul	0.00
20) Perylene-d12	23.31	264	14761	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.96	152	315	0.02	ng/ul	0.00
14) Fluoranthene-d10	19.00	212	821	0.02	ng/ul	0.00
Target Compounds						
15) Fluoranthene	19.03	202	1078	0.022	ng/ul	83
17) Pyrene	19.39	202	1311m	0.020	ng/ul	
18) Benzo(a)anthracene	21.15	228	1144	0.022	ng/ul#	87
19) Chrysene	21.20	228	2623	0.046	ng/ul	95
21) Benzo(b)fluoranthene	22.67	252	2716	0.046	ng/ul#	64
22) Benzo(k)fluoranthene	22.71	252	1335	0.022	ng/ul#	24
23) Benzo(a)pyrene	23.22	252	1436	0.030	ng/ul#	40
24) Indeno(1,2,3-cd)pyrene	25.44	276	1624	0.025	ng/ul#	83
26) Benzo(g,h,i)perylene	26.08	276	1344	0.024	ng/ul#	70

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

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