

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM030620\
 Data File : BM025367.D
 Acq On : 08 Mar 2020 03:02
 Operator : CG/JU
 Sample : L1615-18
 Misc :
 ALS Vial : 53 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 DBDP2

Manual Integrations
 APPROVED

mohammad
 3/9/2020 3:02:35 PM

Quant Time: Mar 08 03:38:38 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 : Sat Mar 07 23:59:42 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.67	152	3641	0.40	ng/ul	0.00
2) Naphthalene-d8	10.43	136	16480	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.29	164	12603	0.40	ng/ul	-0.01
10) Phenanthrene-d10	17.04	188	30378m	0.40	ng/ul	-0.01
16) Chrysene-d12	21.22	240	31237	0.40	ng/ul	0.00
20) Perylene-d12	23.41	264	30908	0.40	ng/ul	-0.01
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.03	152	11159	0.36	ng/ul	0.00
14) Fluoranthene-d10	19.06	212	34382	0.35	ng/ul	-0.01
Target Compounds				Ovalue		
7) Acenaphthylene	14.01	152	1211	0.021	ng/ul#	89
13) Anthracene	17.17	178	4800	0.055	ng/ul	98
15) Fluoranthene	19.10	202	3105	0.025	ng/ul	91
18) Benzo(a)anthracene	21.21	228	2578m	0.022	ng/ul	
19) Chrysene	21.24	228	3367	0.027	ng/ul	95
21) Benzo(b)fluoranthene	22.76	252	11466m	0.093	ng/ul	
22) Benzo(k)fluoranthene	22.80	252	4323m	0.033	ng/ul	
23) Benzo(a)pyrene	23.32	252	4532	0.044	ng/ul#	72
24) Indeno(1,2,3-cd)pyrene	25.58	276	3041	0.023	ng/ul#	100
26) Benzo(g,h,i)perylene	26.25	276	2287	0.021	ng/ul#	87

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

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