

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM040420\
 Data File : BM025934.D
 Acq On : 04 Apr 2020 04:24
 Operator : CG/JU
 Sample : L2062-03
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 F3E02

Manual Integrations
 APPROVED

mohammad
 4/6/2020 1:41:42 PM

Quant Time: Apr 04 05:28:09 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\SOM-EPA-SIM-BM040220.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Apr 04 05:10:01 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.56	152	3173	0.40	ng/ul	0.00
2) Naphthalene-d8	10.32	136	14027	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.20	164	9894	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.94	188	22496m	0.40	ng/ul	0.00
16) Chrysene-d12	21.14	240	25931	0.40	ng/ul	0.00
20) Perylene-d12	23.28	264	28554	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.93	152	9016	0.36	ng/ul	0.00
14) Fluoranthene-d10	18.98	212	25716	0.36	ng/ul	0.00
Target Compounds				Ovalue		
7) Acenaphthylene	13.92	152	948	0.022	ng/ul#	80
15) Fluoranthene	19.00	202	3507	0.040	ng/ul	91
17) Pyrene	19.37	202	7425	0.080	ng/ul#	92
18) Benzo(a)anthracene	21.12	228	4633	0.050	ng/ul	93
19) Chrysene	21.17	228	8328m	0.082	ng/ul	
21) Benzo(b)fluoranthene	22.65	252	16257m	0.155	ng/ul	
22) Benzo(k)fluoranthene	22.69	252	5258m	0.047	ng/ul	
23) Benzo(a)pyrene	23.19	252	6742	0.072	ng/ul#	82
24) Indeno(1,2,3-cd)pyrene	25.40	276	4599	0.037	ng/ul#	100
26) Benzo(g,h,i)perylene	26.04	276	3395	0.032	ng/ul#	88

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

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