

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM031220\
 Data File : BM025531.D
 Acq On : 12 Mar 2020 20:25
 Operator : CG/JU
 Sample : L1782-09
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C0AB3

Manual Integrations
 APPROVED

mohammad
 3/16/2020 1:50:25 PM

Quant Time: Mar 13 05:18:00 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	3581	0.40	ng/ul	0.00
2) Naphthalene-d8	10.42	136	15463	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.28	164	10943	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.03	188	23133m	0.40	ng/ul	0.00
16) Chrysene-d12	21.21	240	16671	0.40	ng/ul	0.00
20) Perylene-d12	23.40	264	17686	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.02	152	9759	0.35	ng/ul	0.00
14) Fluoranthene-d10	19.05	212	25382	0.40	ng/ul	0.00
Target Compounds					Ovalue	
8) Acenaphthene	14.35	153	842	0.021	ng/ul	99
12) Phenanthrene	17.07	178	1703	0.023	ng/ul#	85
13) Anthracene	17.16	178	1854	0.029	ng/ul#	89
15) Fluoranthene	19.08	202	4552	0.057	ng/ul	90
17) Pyrene	19.45	202	3687	0.049	ng/ul#	79
18) Benzo(a)anthracene	21.20	228	2087	0.035	ng/ul#	91
19) Chrysene	21.25	228	2224	0.033	ng/ul#	92
21) Benzo(b)fluoranthene	22.75	252	3367m	0.048	ng/ul	
23) Benzo(a)pyrene	23.30	252	2040	0.035	ng/ul#	42
24) Indeno(1,2,3-cd)pyrene	25.57	276	1661	0.021	ng/ul#	90
26) Benzo(g,h,i)perylene	26.23	276	1559	0.023	ng/ul#	69

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

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