

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM031220\
 Data File : BM025539.D
 Acq On : 13 Mar 2020 01:14
 Operator : CG/JU
 Sample : L1782-08
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C0AB2

Manual Integrations
 APPROVED

mohammad
 3/16/2020 1:50:57 PM

Quant Time: Mar 13 05:34:42 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	3227	0.40	ng/ul	0.00
2) Naphthalene-d8	10.42	136	14332	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.28	164	10831	0.40	ng/ul	-0.01
10) Phenanthrene-d10	17.03	188	24464m	0.40	ng/ul	0.00
16) Chrysene-d12	21.21	240	18230	0.40	ng/ul	0.00
20) Perylene-d12	23.40	264	17431	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.02	152	10727	0.42	ng/ul	0.00
14) Fluoranthene-d10	19.05	212	30635	0.46	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.47	128	5052	0.119	ng/ul	97
5) 2-Methylnaphthalene	12.09	142	1161	0.038	ng/ul	99
7) Acenaphthylene	14.00	152	1718	0.039	ng/ul#	83
8) Acenaphthene	14.35	153	33032	0.851	ng/ul	98
9) Fluorene	15.33	166	26782	0.566	ng/ul	99
12) Phenanthrene	17.07	178	9047	0.117	ng/ul	98
13) Anthracene	17.16	178	6128	0.091	ng/ul	96
15) Fluoranthene	19.08	202	25437	0.303	ng/ul	97
17) Pyrene	19.45	202	14846m	0.181	ng/ul	
18) Benzo(a)anthracene	21.20	228	1694	0.026	ng/ul#	88
19) Chrysene	21.25	228	1800	0.025	ng/ul#	89
21) Benzo(b)fluoranthene	22.75	252	2318m	0.034	ng/ul	
23) Benzo(a)pyrene	23.30	252	1261	0.022	ng/ul#	20

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

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