

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM030920\
 Data File : BM025488.D
 Acq On : 11 Mar 2020 13:14
 Operator : CG/JU
 Sample : L1677-11
 Misc :
 ALS Vial : 81 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C0CC1

Manual Integrations
 APPROVED

mohammad
 3/13/2020 8:48:40 AM

Quant Time: Mar 11 13:56:20 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 : Wed Mar 11 07:15:41 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.66	152	3561	0.40	ng/ul	0.00
2) Naphthalene-d8	10.43	136	15987	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.29	164	11784	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.03	188	25302m	0.40	ng/ul	0.00
16) Chrysene-d12	21.22	240	17934	0.40	ng/ul	0.00
20) Perylene-d12	23.40	264	17228	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.02	152	8502	0.30	ng/ul	0.00
14) Fluoranthene-d10	19.06	212	20454	0.29	ng/ul	0.00
Target Compounds				Ovalue		
7) Acenaphthylene	14.01	152	1001	0.021	ng/ul#	76
12) Phenanthrene	17.07	178	4581	0.057	ng/ul#	95
15) Fluoranthene	19.09	202	16730	0.192	ng/ul	97
17) Pyrene	19.45	202	16044	0.199	ng/ul#	96
18) Benzo(a)anthracene	21.20	228	6758	0.105	ng/ul	95
19) Chrysene	21.26	228	7186	0.100	ng/ul	99
21) Benzo(b)fluoranthene	22.76	252	10264m	0.150	ng/ul	
22) Benzo(k)fluoranthene	22.79	252	4238m	0.059	ng/ul	
23) Benzo(a)pyrene	23.31	252	6340	0.112	ng/ul#	82
24) Indeno(1,2,3-cd)pyrene	25.57	276	4345	0.057	ng/ul#	87
26) Benzo(g,h,i)perylene	26.23	276	4134	0.062	ng/ul#	89

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

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