

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM030920\
 Data File : BM025496.D
 Acq On : 11 Mar 2020 18:05
 Operator : CG/JU
 Sample : L1677-04
 Misc :
 ALS Vial : 89 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C0CC8

Manual Integrations
 APPROVED

mohammad
 3/13/2020 8:49:02 AM

Quant Time: Mar 12 02:06:25 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	3327	0.40	ng/ul	0.00
2) Naphthalene-d8	10.43	136	14078	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.29	164	9904	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.03	188	21374	0.40	ng/ul	0.00
16) Chrysene-d12	21.22	240	15581	0.40	ng/ul	0.00
20) Perylene-d12	23.41	264	15586	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.02	152	6368	0.25	ng/ul	0.00
14) Fluoranthene-d10	19.06	212	15667	0.27	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.48	128	3946	0.095	ng/ul	96
5) 2-Methylnaphthalene	12.09	142	3533	0.119	ng/ul	99
7) Acenaphthylene	14.01	152	2365	0.058	ng/ul#	86
8) Acenaphthene	14.35	153	12173	0.343	ng/ul	99
9) Fluorene	15.34	166	28332	0.655	ng/ul	99
12) Phenanthrene	17.07	178	79729	1.177	ng/ul	100
13) Anthracene	17.16	178	14345	0.243	ng/ul	99
15) Fluoranthene	19.09	202	221383	3.015	ng/ul	100
17) Pyrene	19.45	202	150982	2.159	ng/ul	99
18) Benzo(a)anthracene	21.20	228	30527	0.548	ng/ul	97
19) Chrysene	21.26	228	32987	0.527	ng/ul	98
21) Benzo(b)fluoranthene	22.76	252	34980m	0.567	ng/ul	
22) Benzo(k)fluoranthene	22.79	252	14578m	0.225	ng/ul	
23) Benzo(a)pyrene	23.31	252	22847	0.446	ng/ul	99
24) Indeno(1,2,3-cd)pyrene	25.57	276	16228	0.233	ng/ul#	95
25) Dibenzo(a,h)anthracene	25.58	278	4792	0.082	ng/ul#	79
26) Benzo(a,h,i)perylene	26.24	276	14646	0.243	ng/ul#	97

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

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