

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM030920\
 Data File : BM025460.D
 Acq On : 10 Mar 2020 19:38
 Operator : CG/JU
 Sample : L1617-20
 Misc :
 ALS Vial : 54 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 DBDT4

Manual Integrations
 APPROVED

mohammad
 3/11/2020 8:01:45 AM

Quant Time: Mar 11 00:07:16 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 : Tue Mar 10 23:48:01 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.66	152	3491	0.40	ng/ul	0.00
2) Naphthalene-d8	10.43	136	15307	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.29	164	10230	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.03	188	20927	0.40	ng/ul	0.00
16) Chrysene-d12	21.22	240	16295	0.40	ng/ul	0.00
20) Perylene-d12	23.41	264	19151	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.02	152	3486	0.13	ng/ul	0.00
14) Fluoranthene-d10	19.06	212	7908	0.14	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.48	128	16973	0.375	ng/ul	99
5) 2-Methylnaphthalene	12.10	142	12502	0.386	ng/ul	100
7) Acenaphthylene	14.00	152	59235	1.417	ng/ul	99
8) Acenaphthene	14.35	153	79361	2.165	ng/ul	99
9) Fluorene	15.34	166	108304	2.424	ng/ul	100
12) Phenanthrene	17.07	178	656244	9.895	ng/ul	100
13) Anthracene	17.17	178	1561697	27.059	ng/ul	99
15) Fluoranthene	19.09	202	2819385	39.217	ng/ul	98
17) Pyrene	19.46	202	3275619	44.786	ng/ul	99
18) Benzo(a)anthracene	21.21	228	1490130	25.578	ng/ul	97
19) Chrysene	21.26	228	2027905	30.997	ng/ul	98
21) Benzo(b)fluoranthene	22.76	252	2654271m	35.001	ng/ul	
22) Benzo(k)fluoranthene	22.80	252	855454m	10.766	ng/ul	
23) Benzo(a)pyrene	23.32	252	1180566m	18.747	ng/ul	
24) Indeno(1,2,3-cd)pyrene	25.58	276	778599	9.113	ng/ul#	99
25) Dibenzo(a,h)anthracene	25.59	278	212374	2.955	ng/ul	97
26) Benzo(a,h,i)perylene	26.24	276	507270	6.845	ng/ul	98

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

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