

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM030620\
 Data File : BM025404.D
 Acq On : 09 Mar 2020 02:55
 Operator : CG/JU
 Sample : L1617-21
 Misc :
 ALS Vial : 90 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 DBDT5

Manual Integrations
 APPROVED

mohammad
 3/9/2020 3:04:15 PM

Quant Time: Mar 09 03:58:18 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\SOM-EPA-SIM-BM030320.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Mar 04 08:02:03 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.67	152	3896	0.40	ng/ul	-0.01
2) Naphthalene-d8	10.43	136	17618	0.40	ng/ul	-0.02
6) Acenaphthene-d10	14.29	164	13598	0.40	ng/ul	-0.01
10) Phenanthrene-d10	17.03	188	31343m	0.40	ng/ul	-0.01
16) Chrysene-d12	21.22	240	28061	0.40	ng/ul	-0.01
20) Perylene-d12	23.41	264	24154	0.40	ng/ul	-0.01
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.03	152	10347	0.31	ng/ul	-0.02
14) Fluoranthene-d10	19.06	212	30851	0.30	ng/ul	-0.01
Target Compounds				Ovalue		
3) Naphthalene	10.48	128	225383	4.251	ng/ul#	82
5) 2-Methylnaphthalene	12.10	142	119093	3.084	ng/ul	100
7) Acenaphthylene	14.01	152	10499	0.165	ng/ul	97
8) Acenaphthene	14.35	153	218073	4.416	ng/ul	98
9) Fluorene	15.34	166	185061	2.989	ng/ul	95
12) Phenanthrene	17.08	178	1338194	13.240	ng/ul	95
15) Fluoranthene	19.09	202	537722	4.251	ng/ul	96
17) Pyrene	19.46	202	344635	3.238	ng/ul	96
18) Benzo(a)anthracene	21.20	228	49514	0.480	ng/ul	99
19) Chrysene	21.26	228	31558	0.278	ng/ul	99
21) Benzo(b)fluoranthene	22.76	252	22589m	0.234	ng/ul	
22) Benzo(k)fluoranthene	22.79	252	8013m	0.079	ng/ul	
23) Benzo(a)pyrene	23.31	252	11399	0.140	ng/ul	96
24) Indeno(1,2,3-cd)pyrene	25.58	276	3396	0.034	ng/ul#	100
26) Benzo(g,h,i)perylene	26.24	276	2605	0.030	ng/ul	92

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

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