

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM030920\
 Data File : BM025463.D
 Acq On : 10 Mar 2020 21:28
 Operator : CG/JU
 Sample : L1617-18
 Misc :
 ALS Vial : 57 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 DBDT2

Manual Integrations
 APPROVED

mohammad
 3/11/2020 8:01:52 AM

Quant Time: Mar 11 00:10:35 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	2904	0.40	ng/ul	0.00
2) Naphthalene-d8	10.43	136	12543	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.29	164	8919	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.03	188	19445	0.40	ng/ul	0.00
16) Chrysene-d12	21.22	240	13900	0.40	ng/ul	0.00
20) Perylene-d12	23.40	264	13212	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.03	152	5957	0.26	ng/ul	0.00
14) Fluoranthene-d10	19.06	212	14266	0.27	ng/ul	0.00
Target Compounds				Ovalue		
7) Acenaphthylene	14.01	152	1232	0.034	ng/ul#	88
8) Acenaphthene	14.35	153	2994	0.094	ng/ul	98
9) Fluorene	15.34	166	7893	0.203	ng/ul	100
12) Phenanthrene	17.07	178	70711	1.147	ng/ul	99
13) Anthracene	17.16	178	13721	0.256	ng/ul	98
15) Fluoranthene	19.09	202	57850	0.866	ng/ul	99
17) Pyrene	19.45	202	46891	0.752	ng/ul	98
18) Benzo(a)anthracene	21.20	228	13695	0.276	ng/ul	97
19) Chrysene	21.25	228	18972	0.340	ng/ul	99
21) Benzo(b)fluoranthene	22.75	252	39000m	0.745	ng/ul	
22) Benzo(k)fluoranthene	22.79	252	17048m	0.311	ng/ul	
23) Benzo(a)pyrene	23.31	252	12092	0.278	ng/ul	96
24) Indeno(1,2,3-cd)pyrene	25.57	276	14046	0.238	ng/ul#	98
25) Dibenzo(a,h)anthracene	25.58	278	3718	0.075	ng/ul#	79
26) Benzo(g,h,i)perylene	26.23	276	7631	0.149	ng/ul#	94

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

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