

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM030920\
 Data File : BM025444.D
 Acq On : 10 Mar 2020 09:54
 Operator : CG/JU
 Sample : L1612-17
 Misc :
 ALS Vial : 38 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 DBDL9

Manual Integrations
 APPROVED

mohammad
 3/11/2020 8:01:06 AM

Quant Time: Mar 10 12:13:39 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	3077	0.40	ng/ul	0.00
2) Naphthalene-d8	10.43	136	12889	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.29	164	8482	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.03	188	16568	0.40	ng/ul	0.00
16) Chrysene-d12	21.22	240	12445	0.40	ng/ul	0.00
20) Perylene-d12	23.40	264	13041	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.02	152	7226	0.31	ng/ul	0.00
14) Fluoranthene-d10	19.06	212	14559	0.32	ng/ul	0.00
Target Compounds				Ovalue		
7) Acenaphthylene	14.01	152	2579	0.074	ng/ul#	92
9) Fluorene	15.34	166	741	0.020	ng/ul#	96
12) Phenanthrene	17.07	178	4032	0.077	ng/ul	96
13) Anthracene	17.16	178	10017	0.219	ng/ul	98
15) Fluoranthene	19.09	202	20059	0.352	ng/ul	98
17) Pyrene	19.45	202	32550	0.583	ng/ul	98
18) Benzo(a)anthracene	21.20	228	16330	0.367	ng/ul	97
19) Chrysene	21.26	228	12904	0.258	ng/ul	98
21) Benzo(b)fluoranthene	22.76	252	51400m	0.995	ng/ul	
22) Benzo(k)fluoranthene	22.80	252	18763m	0.347	ng/ul	
23) Benzo(a)pyrene	23.31	252	17540	0.409	ng/ul	99
24) Indeno(1,2,3-cd)pyrene	25.57	276	13270	0.228	ng/ul#	99
25) Dibenzo(a,h)anthracene	25.58	278	3841	0.078	ng/ul#	82
26) Benzo(g,h,i)perylene	26.23	276	7458	0.148	ng/ul#	95

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

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