

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM030320\
 Data File : BM025278.D
 Acq On : 05 Mar 2020 08:35
 Operator : CG/JU
 Sample : L1543-12
 Misc :
 ALS Vial : 74 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 DBD94

Manual Integrations
 APPROVED

mohammad
 3/5/2020 3:31:01 PM

Quant Time: Mar 05 14:06:10 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	3428	0.40	ng/ul	0.00
2) Naphthalene-d8	10.44	136	14929	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.30	164	11280	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.04	188	27284m	0.40	ng/ul	0.00
16) Chrysene-d12	21.23	240	26817	0.40	ng/ul	0.00
20) Perylene-d12	23.42	264	24996	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.04	152	7184	0.26	ng/ul	0.00
14) Fluoranthene-d10	19.07	212	25892	0.29	ng/ul	0.00
Target Compounds				Ovalue		
7) Acenaphthylene	14.02	152	1619	0.031	ng/ul	94
12) Phenanthrene	17.08	178	2313	0.026	ng/ul	95
13) Anthracene	17.17	178	6008	0.077	ng/ul	97
15) Fluoranthene	19.10	202	32594	0.296	ng/ul	96
17) Pyrene	19.46	202	49192m	0.484	ng/ul	
18) Benzo(a)anthracene	21.21	228	22030	0.223	ng/ul	98
19) Chrysene	21.27	228	38123	0.352	ng/ul	99
21) Benzo(b)fluoranthene	22.77	252	75651m	0.756	ng/ul	
22) Benzo(k)fluoranthene	22.81	252	25291m	0.241	ng/ul	
23) Benzo(a)pyrene	23.33	252	21978	0.262	ng/ul#	94
24) Indeno(1,2,3-cd)pyrene	25.59	276	18992	0.181	ng/ul#	100
25) Dibenzo(a,h)anthracene	25.60	278	4523	0.052	ng/ul#	79
26) Benzo(g,h,i)perylene	26.26	276	11803	0.133	ng/ul#	95

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

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