

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM031720\
 Data File : BM025614.D
 Acq On : 17 Mar 2020 17:20
 Operator : CG/JU
 Sample : L1516-04
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 DBD22

Manual Integrations
 APPROVED

mohammad
 3/18/2020 9:01:54 AM

Quant Time: Mar 17 17:55:48 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 17 17:05:05 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.59	152	2063	0.40	ng/ul	0.00
2) Naphthalene-d8	10.35	136	7942	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.22	164	5245	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.97	188	12250m	0.40	ng/ul	0.00
16) Chrysene-d12	21.16	240	13429	0.40	ng/ul	0.00
20) Perylene-d12	23.31	264	13999	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.95	152	2943	0.21	ng/ul	0.00
14) Fluoranthene-d10	19.00	212	7608	0.23	ng/ul	0.00
Target Compounds				Ovalue		
5) 2-Methylnaphthalene	12.03	142	896	0.053	ng/ul	98
7) Acenaphthylene	13.94	152	2538	0.118	ng/ul#	84
8) Acenaphthene	14.28	153	52402	2.788	ng/ul	99
9) Fluorene	15.27	166	38450	1.679	ng/ul	99
12) Phenanthrene	17.01	178	147607	3.802	ng/ul	99
13) Anthracene	17.10	178	15780	0.467	ng/ul	99
15) Fluoranthene	19.03	202	235071	5.586	ng/ul	99
17) Pyrene	19.39	202	147530	2.448	ng/ul	99
18) Benzo(a)anthracene	21.14	228	22495	0.469	ng/ul	100
19) Chrysene	21.20	228	20294	0.376	ng/ul	98
21) Benzo(b)fluoranthene	22.68	252	9231m	0.167	ng/ul	
22) Benzo(k)fluoranthene	22.72	252	4125m	0.071	ng/ul	
23) Benzo(a)pyrene	23.22	252	3414	0.074	ng/ul#	72
24) Indeno(1,2,3-cd)pyrene	25.44	276	1595	0.026	ng/ul#	76
26) Benzo(g,h,i)perylene	26.09	276	1220	0.023	ng/ul#	69

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

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