

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM030620\
 Data File : BM025332.D
 Acq On : 07 Mar 2020 03:54
 Operator : CG/JU
 Sample : L1612-18
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 DBDM0

Manual Integrations
 APPROVED

mohammad
 3/9/2020 3:00:42 PM

Quant Time: Mar 07 06:53:36 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 : Sat Mar 07 00:01:42 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.67	152	3778	0.40	ng/ul	0.00
2) Naphthalene-d8	10.44	136	16759	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.30	164	12215	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.04	188	29226	0.40	ng/ul	0.00
16) Chrysene-d12	21.22	240	28360	0.40	ng/ul	0.00
20) Perylene-d12	23.41	264	25464	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.03	152	3277	0.10	ng/ul	0.00
14) Fluoranthene-d10	19.07	212	10063	0.11	ng/ul	0.00
Target Compounds				Ovalue		
5) 2-Methylnaphthalene	12.11	142	1643	0.045	ng/ul	96
7) Acenaphthylene	14.01	152	2657	0.046	ng/ul	99
9) Fluorene	15.35	166	4283	0.077	ng/ul	96
12) Phenanthrene	17.08	178	14761	0.157	ng/ul	97
13) Anthracene	17.17	178	86089	1.029	ng/ul	95
15) Fluoranthene	19.10	202	15281	0.130	ng/ul	97
17) Pyrene	19.46	202	31253m	0.291	ng/ul	
18) Benzo(a)anthracene	21.21	228	87188	0.836	ng/ul	100
19) Chrysene	21.26	228	112800m	0.985	ng/ul	
21) Benzo(b)fluoranthene	22.76	252	233624m	2.293	ng/ul	
22) Benzo(k)fluoranthene	22.80	252	91623m	0.858	ng/ul	
23) Benzo(a)pyrene	23.32	252	68791	0.804	ng/ul#	90
24) Indeno(1,2,3-cd)pyrene	25.58	276	47666	0.447	ng/ul#	100
25) Dibenzo(a,h)anthracene	25.59	278	13787	0.156	ng/ul	96
26) Benzo(g,h,i)perylene	26.25	276	23290	0.257	ng/ul#	92

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

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