

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM031820\
 Data File : BM025646.D
 Acq On : 19 Mar 2020 01:58
 Operator : CG/JU
 Sample : L1786-12DL 5X
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C0AC8DL

Manual Integrations
 APPROVED

mohammad
 3/19/2020 12:44:00 PM

Quant Time: Mar 19 02:50:05 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 : Thu Mar 19 00:21:10 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.58	152	2374	0.40	ng/ul	0.00
2) Naphthalene-d8	10.35	136	9723	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.22	164	6308	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.96	188	14722	0.40	ng/ul	0.00
16) Chrysene-d12	21.15	240	15651	0.40	ng/ul	0.00
20) Perylene-d12	23.30	264	16103	0.40	ng/ul	-0.01
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.95	152	568	0.03	ng/ul	0.00
14) Fluoranthene-d10	19.00	212	1534	0.04	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.40	128	1186	0.041	ng/ul#	86
5) 2-Methylnaphthalene	12.02	142	460	0.022	ng/ul	94
7) Acenaphthylene	13.94	152	2506	0.097	ng/ul#	84
8) Acenaphthene	14.28	153	1221	0.054	ng/ul	97
9) Fluorene	15.27	166	3629	0.132	ng/ul	98
12) Phenanthrene	17.00	178	43237	0.927	ng/ul	100
13) Anthracene	17.09	178	24907	0.613	ng/ul	99
15) Fluoranthene	19.02	202	102166	2.020	ng/ul	98
17) Pyrene	19.38	202	78297	1.115	ng/ul	98
18) Benzo(a)anthracene	21.14	228	44730	0.799	ng/ul	99
19) Chrysene	21.19	228	47745	0.760	ng/ul	98
21) Benzo(b)fluoranthene	22.67	252	46164m	0.724	ng/ul	
22) Benzo(k)fluoranthene	22.70	252	21547m	0.323	ng/ul	
23) Benzo(a)pyrene	23.21	252	28243	0.533	ng/ul	96
24) Indeno(1,2,3-cd)pyrene	25.42	276	18427	0.257	ng/ul#	99
25) Dibenzo(a,h)anthracene	25.43	278	6015	0.100	ng/ul#	87
26) Benzo(a,h,i)perylene	26.07	276	11759	0.189	ng/ul#	96

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

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