

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM031820\
 Data File : BM025639.D
 Acq On : 18 Mar 2020 21:46
 Operator : CG/JU
 Sample : L1786-13DL 5X
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C0AC9DL

Manual Integrations
 APPROVED

mohammad
 3/19/2020 12:43:38 PM

Quant Time: Mar 19 00:32:15 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 : Wed Mar 18 05:07:34 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.58	152	2507	0.40	ng/ul	0.00
2) Naphthalene-d8	10.35	136	10041	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.22	164	6388	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.96	188	15000	0.40	ng/ul	0.00
16) Chrysene-d12	21.15	240	16341	0.40	ng/ul	0.00
20) Perylene-d12	23.30	264	17797	0.40	ng/ul	-0.01
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.95	152	365	0.02	ng/ul	0.00
14) Fluoranthene-d10	19.00	212	1011	0.02	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.40	128	7261	0.244	ng/ul	100
5) 2-Methylnaphthalene	12.02	142	2160	0.102	ng/ul	98
7) Acenaphthylene	13.94	152	6651	0.255	ng/ul#	86
8) Acenaphthene	14.28	153	8021	0.350	ng/ul	99
9) Fluorene	15.27	166	13747	0.493	ng/ul	100
12) Phenanthrene	17.00	178	187299	3.940	ng/ul	100
13) Anthracene	17.09	178	52480	1.269	ng/ul	99
15) Fluoranthene	19.03	202	406560	7.890	ng/ul	99
17) Pyrene	19.39	202	325970	4.444	ng/ul	99
18) Benzo(a)anthracene	21.14	228	209476	3.585	ng/ul	99
19) Chrysene	21.19	228	193538	2.950	ng/ul	97
21) Benzo(b)fluoranthene	22.67	252	259651m	3.684	ng/ul	
22) Benzo(k)fluoranthene	22.71	252	94340m	1.278	ng/ul	
23) Benzo(a)pyrene	23.21	252	177339m	3.030	ng/ul	
24) Indeno(1,2,3-cd)pyrene	25.43	276	119271	1.502	ng/ul#	99
25) Dibenzo(a,h)anthracene	25.43	278	34121	0.511	ng/ul	98
26) Benzo(a,h,i)perylene	26.08	276	85611	1.243	ng/ul	98

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

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