

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN102419\
 Data File : BN008418.D
 Acq On : 25 Oct 2019 09:18
 Operator : JU
 Sample : K5236-11DL 5X
 Misc :
 ALS Vial : 40 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 BEPM1DL

Manual Integrations
 APPROVED

mohammad
 10/25/2019 9:28:31 PM

Quant Time: Oct 25 12:12:23 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN101019.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Oct 25 00:45:13 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.57	152	2134	0.40	ng/ul	0.00
2) Naphthalene-d8	10.33	136	10839	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.21	164	7107	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.96	188	15099	0.40	ng/ul	0.00
16) Chrysene-d12	21.17	240	12403	0.40	ng/ul	0.00
20) Perylene-d12	23.34	264	9830	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.94	152	559	0.03	ng/ul	0.00
14) Fluoranthene-d10	19.00	212	1899	0.04	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.38	128	1662	0.054	ng/ul#	88
5) 2-Methylnaphthalene	12.02	142	1423	0.067	ng/ul	100
7) Acenaphthylene	13.93	152	8385	0.248	ng/ul	99
8) Acenaphthene	14.27	153	2793	0.106	ng/ul	97
9) Fluorene	15.26	166	5911	0.198	ng/ul	97
12) Phenanthrene	17.00	178	101192	2.289	ng/ul	99
13) Anthracene	17.09	178	18269	0.470	ng/ul	99
15) Fluoranthene	19.03	202	198711	3.403	ng/ul	100
17) Pyrene	19.39	202	154650	3.132	ng/ul	100
18) Benzo(a)anthracene	21.15	228	80724	1.790	ng/ul	98
19) Chrysene	21.20	228	80244	1.450	ng/ul	98
21) Benzo(b)fluoranthene	22.69	252	90451m	2.827	ng/ul	
22) Benzo(k)fluoranthene	22.73	252	33483m	0.924	ng/ul	
23) Benzo(a)pyrene	23.24	252	56459m	1.677	ng/ul	
24) Indeno(1,2,3-cd)pyrene	25.49	276	35327	0.891	ng/ul	99
25) Dibenzo(a,h)anthracene	25.49	278	8897	0.284	ng/ul#	85
26) Benzo(a,h,i)perylene	26.14	276	32626	0.897	ng/ul	98

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

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