

Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN102419\
 Data File : BN008389.D
 Acq On : 24 Oct 2019 16:01
 Operator : JU
 Sample : K5235-14DL 5X
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 BEPB4DL

Manual Integrations
 APPROVED

mohammad
 10/25/2019 8:05:46 AM

Quant Time: Oct 24 16:35:07 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 : Thu Oct 24 13:52:48 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.58	152	2178	0.40	ng/ul	0.00
2) Naphthalene-d8	10.34	136	10470	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.21	164	6501	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.96	188	13769	0.40	ng/ul	0.00
16) Chrysene-d12	21.17	240	11195	0.40	ng/ul	0.00
20) Perylene-d12	23.34	264	9196	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.95	152	682	0.04	ng/ul	0.00
14) Fluoranthene-d10	19.00	212	2035	0.04	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	10.39	128	1780	0.060	ng/ul#	90
5) 2-Methylnaphthalene	12.02	142	929	0.045	ng/ul	100
7) Acenaphthylene	13.93	152	5306	0.171	ng/ul	98
8) Acenaphthene	14.27	153	2345	0.097	ng/ul	98
9) Fluorene	15.26	166	3152	0.115	ng/ul#	96
12) Phenanthrene	17.00	178	67636	1.678	ng/ul	99
13) Anthracene	17.09	178	14151	0.399	ng/ul	99
15) Fluoranthene	19.03	202	182873	3.435	ng/ul	98
17) Pyrene	19.39	202	199540	4.477	ng/ul	98
18) Benzo(a)anthracene	21.15	228	104874	2.577	ng/ul	98
19) Chrysene	21.20	228	113781	2.278	ng/ul	98
21) Benzo(b)fluoranthene	22.69	252	109210m	3.648	ng/ul	
22) Benzo(k)fluoranthene	22.73	252	39741m	1.172	ng/ul	
23) Benzo(a)pyrene	23.24	252	78700m	2.498	ng/ul	
24) Indeno(1,2,3-cd)pyrene	25.48	276	43634	1.176	ng/ul#	99
25) Dibenzo(a,h)anthracene	25.49	278	12131	0.414	ng/ul	94
26) Benzo(g,h,i)perylene	26.14	276	45410	1.335	ng/ul	98

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

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