

Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN102419\
 Data File : BN008396.D
 Acq On : 24 Oct 2019 20:13
 Operator : JU
 Sample : K5235-18DL 10X
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 BEPB8DL

Manual Integrations
 APPROVED

mohammad
 10/25/2019 8:06:03 AM

Quant Time: Oct 24 22:08:55 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 : Wed Oct 23 14:07:55 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.58	152	1583	0.40	ng/ul	-0.02
2) Naphthalene-d8	10.34	136	7721	0.40	ng/ul	-0.02
6) Acenaphthene-d10	14.21	164	5010	0.40	ng/ul	-0.01
10) Phenanthrene-d10	16.96	188	10478	0.40	ng/ul	-0.02
16) Chrysene-d12	21.17	240	9523	0.40	ng/ul	-0.01
20) Perylene-d12	23.33	264	7801	0.40	ng/ul	-0.02
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.95	152	280	0.02	ng/ul	0.00
14) Fluoranthene-d10	19.00	212	935	0.03	ng/ul	-0.02
Target Compounds				Qvalue		
3) Naphthalene	10.39	128	608	0.028	ng/ul#	67
5) 2-Methylnaphthalene	12.03	142	372	0.025	ng/ul	96
7) Acenaphthylene	13.93	152	1581	0.066	ng/ul#	89
8) Acenaphthene	14.27	153	895	0.048	ng/ul	96
9) Fluorene	15.27	166	1039	0.049	ng/ul#	94
12) Phenanthrene	17.00	178	17808	0.580	ng/ul	99
13) Anthracene	17.09	178	4500	0.167	ng/ul	96
15) Fluoranthene	19.02	202	37953	0.937	ng/ul	96
17) Pyrene	19.39	202	42214	1.113	ng/ul	100
18) Benzo(a)anthracene	21.15	228	20723	0.599	ng/ul	97
19) Chrysene	21.20	228	23278	0.548	ng/ul	98
21) Benzo(b)fluoranthene	22.69	252	23006m	0.906	ng/ul	
22) Benzo(k)fluoranthene	22.73	252	6625m	0.230	ng/ul	
23) Benzo(a)pyrene	23.24	252	15799	0.591	ng/ul	96
24) Indeno(1,2,3-cd)pyrene	25.48	276	9435	0.300	ng/ul#	99
25) Dibenzo(a,h)anthracene	25.49	278	2282	0.092	ng/ul#	75
26) Benzo(g,h,i)perylene	26.14	276	9558	0.331	ng/ul	96

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

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