

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN102419\
 Data File : BN008406.D
 Acq On : 25 Oct 2019 02:10
 Operator : JU
 Sample : K5236-04DL 5X
 Misc :
 ALS Vial : 29 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 BEPL3DL

Manual Integrations
 APPROVED

mohammad
 10/25/2019 8:06:29 AM

Quant Time: Oct 25 03:07:59 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	7.57	152	1998	0.40	ng/ul	-0.02
2) Naphthalene-d8	10.33	136	9854	0.40	ng/ul	-0.02
6) Acenaphthene-d10	14.21	164	6479	0.40	ng/ul	-0.01
10) Phenanthrene-d10	16.96	188	13750	0.40	ng/ul	-0.02
16) Chrysene-d12	21.16	240	11679	0.40	ng/ul	-0.02
20) Perylene-d12	23.33	264	9822	0.40	ng/ul	-0.02
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.94	152	667	0.04	ng/ul	-0.02
14) Fluoranthene-d10	19.00	212	1940	0.04	ng/ul	-0.02
Target Compounds				Ovalue		
3) Naphthalene	10.38	128	4643	0.166	ng/ul	97
5) 2-Methylnaphthalene	12.01	142	2804	0.146	ng/ul	100
7) Acenaphthylene	13.93	152	6213	0.201	ng/ul	97
8) Acenaphthene	14.27	153	3264	0.136	ng/ul	97
9) Fluorene	15.26	166	3346	0.123	ng/ul#	95
12) Phenanthrene	17.00	178	58775	1.460	ng/ul	100
13) Anthracene	17.09	178	11570	0.327	ng/ul	95
15) Fluoranthene	19.03	202	97504	1.834	ng/ul	100
17) Pyrene	19.39	202	81339	1.749	ng/ul	100
18) Benzo(a)anthracene	21.15	228	40856	0.962	ng/ul	93
19) Chrysene	21.20	228	41798m	0.802	ng/ul	
21) Benzo(b)fluoranthene	22.69	252	54903m	1.717	ng/ul	
22) Benzo(k)fluoranthene	22.73	252	14830m	0.409	ng/ul	
23) Benzo(a)pyrene	23.24	252	34340	1.021	ng/ul#	88
24) Indeno(1,2,3-cd)pyrene	25.48	276	28152	0.710	ng/ul	98
25) Dibenzo(a,h)anthracene	25.48	278	6083	0.194	ng/ul#	49
26) Benzo(a,h,i)perylene	26.14	276	28696	0.790	ng/ul#	94

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

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