

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN102419\
 Data File : BN008410.D
 Acq On : 25 Oct 2019 04:33
 Operator : JU
 Sample : K5236-05DL 5X
 Misc :
 ALS Vial : 32 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 BEPL5DL

Manual Integrations
 APPROVED

mohammad
 10/25/2019 9:28:05 PM

Quant Time: Oct 25 11:51:37 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	7.58	152	1922	0.40	ng/ul	0.00
2) Naphthalene-d8	10.33	136	9618	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.21	164	6395	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.96	188	13791	0.40	ng/ul	0.00
16) Chrysene-d12	21.17	240	11787	0.40	ng/ul	0.00
20) Perylene-d12	23.33	264	9496	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.94	152	495	0.03	ng/ul	0.00
14) Fluoranthene-d10	19.00	212	1604	0.03	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.38	128	1101	0.040	ng/ul#	83
5) 2-Methylnaphthalene	12.01	142	735	0.039	ng/ul	98
7) Acenaphthylene	13.93	152	4359	0.143	ng/ul	96
8) Acenaphthene	14.27	153	630	0.027	ng/ul#	94
9) Fluorene	15.26	166	1031	0.038	ng/ul#	92
12) Phenanthrene	17.00	178	25613	0.634	ng/ul	100
13) Anthracene	17.09	178	4946	0.139	ng/ul#	94
15) Fluoranthene	19.02	202	60773	1.140	ng/ul	96
17) Pyrene	19.39	202	49876	1.063	ng/ul	100
18) Benzo(a)anthracene	21.15	228	24641	0.575	ng/ul	94
19) Chrysene	21.20	228	26612m	0.506	ng/ul	
21) Benzo(b)fluoranthene	22.69	252	32018m	1.036	ng/ul	
22) Benzo(k)fluoranthene	22.73	252	11507m	0.329	ng/ul	
23) Benzo(a)pyrene	23.24	252	20120	0.619	ng/ul#	95
24) Indeno(1,2,3-cd)pyrene	25.48	276	15367	0.401	ng/ul#	97
25) Dibenzo(a,h)anthracene	25.49	278	3326	0.110	ng/ul#	54
26) Benzo(a,h,i)perylene	26.14	276	16036	0.456	ng/ul#	94

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

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