

Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN102219\
 Data File : BN008358.D
 Acq On : 23 Oct 2019 20:20
 Operator : JU
 Sample : K5223-19DL 2X
 Misc :
 ALS Vial : 51 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 BEPA7DL

Manual Integrations
 APPROVED

mohammad
 10/24/2019 8:06:53 AM

Quant Time: Oct 24 02:00:27 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	1864	0.40	ng/ul	-0.01
2) Naphthalene-d8	10.34	136	9082	0.40	ng/ul	-0.02
6) Acenaphthene-d10	14.21	164	5825	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.96	188	12662	0.40	ng/ul	-0.01
16) Chrysene-d12	21.17	240	11129	0.40	ng/ul	-0.01
20) Perylene-d12	23.34	264	9028	0.40	ng/ul	-0.01
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.95	152	1686	0.11	ng/ul	-0.01
14) Fluoranthene-d10	19.00	212	5107	0.12	ng/ul	-0.01
Target Compounds					Qvalue	
3) Naphthalene	10.39	128	539	0.021	ng/ul#	68
7) Acenaphthylene	13.93	152	1735	0.063	ng/ul#	91
8) Acenaphthene	14.27	153	573	0.027	ng/ul	95
9) Fluorene	15.27	166	759	0.031	ng/ul#	91
12) Phenanthrene	17.00	178	18199	0.491	ng/ul	99
13) Anthracene	17.09	178	3280	0.101	ng/ul	94
15) Fluoranthene	19.03	202	39066	0.798	ng/ul	96
17) Pyrene	19.39	202	44428	1.003	ng/ul	100
18) Benzo(a)anthracene	21.15	228	23751	0.587	ng/ul	97
19) Chrysene	21.21	228	26409	0.532	ng/ul	98
21) Benzo(b)fluoranthene	22.70	252	23791m	0.810	ng/ul	
22) Benzo(k)fluoranthene	22.73	252	7720m	0.232	ng/ul	
23) Benzo(a)pyrene	23.24	252	17014	0.550	ng/ul	95
24) Indeno(1,2,3-cd)pyrene	25.49	276	9258	0.254	ng/ul#	93
25) Dibenzo(a,h)anthracene	25.49	278	2343	0.081	ng/ul#	75
26) Benzo(g,h,i)perylene	26.14	276	7308	0.219	ng/ul#	94

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

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