

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN102219\
 Data File : BN008342.D
 Acq On : 23 Oct 2019 08:07
 Operator : JU
 Sample : K5223-17
 Misc :
 ALS Vial : 36 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 BEPA5

Manual Integrations
 APPROVED

mohammad
 10/24/2019 8:06:18 AM

Quant Time: Oct 23 09:36:49 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 : Mon Oct 21 23:58:15 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.58	152	1706	0.40	ng/ul	-0.01
2) Naphthalene-d8	10.35	136	8597	0.40	ng/ul	-0.01
6) Acenaphthene-d10	14.22	164	5411	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.97	188	11889m	0.40	ng/ul	0.00
16) Chrysene-d12	21.17	240	11070	0.40	ng/ul	0.00
20) Perylene-d12	23.34	264	9535	0.40	ng/ul	-0.01
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.95	152	3504	0.24	ng/ul	-0.01
14) Fluoranthene-d10	19.00	212	10298	0.25	ng/ul	-0.01
Target Compounds				Ovalue		
3) Naphthalene	10.39	128	1195	0.049	ng/ul#	84
5) 2-Methylnaphthalene	12.02	142	450	0.027	ng/ul	96
7) Acenaphthylene	13.93	152	1879	0.073	ng/ul#	91
8) Acenaphthene	14.28	153	1339	0.067	ng/ul	99
9) Fluorene	15.27	166	1456	0.064	ng/ul#	94
12) Phenanthrene	17.00	178	28883	0.830	ng/ul	99
13) Anthracene	17.10	178	5879	0.192	ng/ul	97
15) Fluoranthene	19.03	202	69051	1.502	ng/ul	99
17) Pyrene	19.39	202	77176	1.751	ng/ul	97
18) Benzo(a)anthracene	21.15	228	42515	1.056	ng/ul	98
19) Chrysene	21.21	228	48274	0.977	ng/ul	99
21) Benzo(b)fluoranthene	22.70	252	46553m	1.500	ng/ul	
22) Benzo(k)fluoranthene	22.74	252	13789m	0.392	ng/ul	
23) Benzo(a)pyrene	23.25	252	32894	1.007	ng/ul	95
24) Indeno(1,2,3-cd)pyrene	25.50	276	18043	0.469	ng/ul	96
25) Dibenzo(a,h)anthracene	25.50	278	4551	0.150	ng/ul#	82
26) Benzo(a,h,i)perylene	26.15	276	14233	0.403	ng/ul	97

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

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