

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN123019\
 Data File : BN009257.D
 Acq On : 30 Dec 2019 13:31
 Operator : JU
 Sample : K6322-05
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 C0C35

Manual Integrations
 APPROVED

mohammad
 12/31/2019 8:27:52 AM

Quant Time: Dec 30 15:24:09 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN121619.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Dec 30 11:05:47 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.51	152	2220	0.40	ng/ul	0.00
2) Naphthalene-d8	10.26	136	8016	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.14	164	4216	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.89	188	8420	0.40	ng/ul	0.00
16) Chrysene-d12	21.09	240	6583	0.40	ng/ul	0.00
20) Perylene-d12	23.22	264	6720	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.86	152	1866	0.13	ng/ul	0.00
14) Fluoranthene-d10	18.92	212	3989	0.16	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.31	128	8527	0.351	ng/ul	98
5) 2-Methylnaphthalene	11.93	142	3536	0.192	ng/ul	100
7) Acenaphthylene	13.85	152	3740	0.169	ng/ul	95
9) Fluorene	15.19	166	613	0.030	ng/ul#	90
12) Phenanthrene	16.93	178	8784	0.300	ng/ul	99
13) Anthracene	17.02	178	12241	0.466	ng/ul	99
15) Fluoranthene	18.95	202	5362	0.160	ng/ul	94
17) Pyrene	19.32	202	7373	0.220	ng/ul	98
18) Benzo(a)anthracene	21.07	228	7199	0.246	ng/ul#	83
19) Chrysene	21.12	228	5133m	0.175	ng/ul	
21) Benzo(b)fluoranthene	22.59	252	57392m	1.829	ng/ul	
22) Benzo(k)fluoranthene	22.63	252	22269m	0.704	ng/ul	
23) Benzo(a)pyrene	23.12	252	25875	0.941	ng/ul	97
24) Indeno(1,2,3-cd)pyrene	25.31	276	15638	0.522	ng/ul#	90
25) Dibenzo(a,h)anthracene	25.31	278	4271	0.179	ng/ul#	80
26) Benzo(a,h,i)perylene	25.95	276	11390	0.452	ng/ul	98

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

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