

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN102219\
 Data File : BN008308.D
 Acq On : 22 Oct 2019 10:39
 Operator : JU
 Sample : K5199-07DL 5X
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 BEP74DL

Manual Integrations
 APPROVED

mohammad
 10/23/2019 8:06:19 AM

Quant Time: Oct 22 13:41:36 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 : Tue Oct 22 13:04:25 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.58	152	2006	0.40	ng/ul	0.00
2) Naphthalene-d8	10.35	136	9694	0.40	ng/ul	-0.01
6) Acenaphthene-d10	14.22	164	6342	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.97	188	13903	0.40	ng/ul	0.00
16) Chrysene-d12	21.18	240	13535	0.40	ng/ul	0.00
20) Perylene-d12	23.35	264	13812	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.95	152	976	0.06	ng/ul	0.00
14) Fluoranthene-d10	19.00	212	3252	0.07	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.39	128	1520	0.055	ng/ul#	86
5) 2-Methylnaphthalene	12.03	142	1208	0.064	ng/ul	98
7) Acenaphthylene	13.94	152	785	0.026	ng/ul#	74
8) Acenaphthene	14.28	153	2497	0.106	ng/ul	99
9) Fluorene	15.27	166	1985	0.074	ng/ul#	96
12) Phenanthrene	17.01	178	24944	0.613	ng/ul	99
13) Anthracene	17.10	178	6048	0.169	ng/ul	98
15) Fluoranthene	19.04	202	26709	0.497	ng/ul	98
17) Pyrene	19.40	202	30991	0.575	ng/ul	99
18) Benzo(a)anthracene	21.16	228	14462	0.294	ng/ul	96
19) Chrysene	21.21	228	16418	0.272	ng/ul	98
21) Benzo(b)fluoranthene	22.71	252	14947m	0.332	ng/ul	
22) Benzo(k)fluoranthene	22.74	252	5311m	0.104	ng/ul	
23) Benzo(a)pyrene	23.26	252	10785	0.228	ng/ul#	92
24) Indeno(1,2,3-cd)pyrene	25.51	276	5790	0.104	ng/ul#	92
25) Dibenzo(a,h)anthracene	25.51	278	1489	0.034	ng/ul#	53
26) Benzo(a,h,i)perylene	26.16	276	5150	0.101	ng/ul#	91

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

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