

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN102219\
 Data File : BN008319.D
 Acq On : 22 Oct 2019 17:46
 Operator : JU
 Sample : K5199-20DL 2X
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 BEPH0DL

Manual Integrations
 APPROVED

mohammad
 10/24/2019 8:05:15 AM

Quant Time: Oct 23 01:46:26 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	2316	0.40	ng/ul	-0.01
2) Naphthalene-d8	10.35	136	11345	0.40	ng/ul	-0.01
6) Acenaphthene-d10	14.22	164	7251	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.97	188	16198	0.40	ng/ul	0.00
16) Chrysene-d12	21.18	240	16297	0.40	ng/ul	0.00
20) Perylene-d12	23.35	264	16130	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.95	152	2234	0.12	ng/ul	-0.01
14) Fluoranthene-d10	19.00	212	7559	0.14	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.39	128	3124	0.097	ng/ul	95
5) 2-Methylnaphthalene	12.02	142	1480	0.067	ng/ul	99
7) Acenaphthylene	13.94	152	9964	0.289	ng/ul	97
8) Acenaphthene	14.28	153	4658	0.173	ng/ul	98
9) Fluorene	15.27	166	6313	0.207	ng/ul	97
12) Phenanthrene	17.01	178	132663	2.797	ng/ul	99
13) Anthracene	17.10	178	26687	0.639	ng/ul	99
15) Fluoranthene	19.04	202	311907	4.980	ng/ul	100
17) Pyrene	19.40	202	290052	4.470	ng/ul	98
18) Benzo(a)anthracene	21.16	228	175574	2.963	ng/ul	97
19) Chrysene	21.21	228	185819	2.556	ng/ul	99
21) Benzo(b)fluoranthene	22.71	252	225782m	4.300	ng/ul	
22) Benzo(k)fluoranthene	22.74	252	73145m	1.230	ng/ul	
23) Benzo(a)pyrene	23.26	252	155818	2.820	ng/ul	91
24) Indeno(1,2,3-cd)pyrene	25.50	276	92299	1.418	ng/ul#	99
25) Dibenzo(a,h)anthracene	25.51	278	24418	0.475	ng/ul	96
26) Benzo(a,h,i)perylene	26.16	276	69675	1.168	ng/ul	98

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

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