

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN101119\
 Data File : BN008113.D
 Acq On : 12 Oct 2019 15:21
 Operator : JU
 Sample : K5178-04
 Misc :
 ALS Vial : 46 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 BEPE7

Manual Integrations
 APPROVED

mohammad
 10/15/2019 7:58:48 AM

Quant Time: Oct 13 03:39:05 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 : Fri Oct 11 19:05:54 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.63	152	2555	0.40	ng/ul	0.00
2) Naphthalene-d8	10.39	136	12395	0.40	ng/ul	-0.01
6) Acenaphthene-d10	14.26	164	8194	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.01	188	18808m	0.40	ng/ul	0.00
16) Chrysene-d12	21.21	240	19853m	0.40	ng/ul	0.00
20) Perylene-d12	23.40	264	18742	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.00	152	3944	0.19	ng/ul	0.00
14) Fluoranthene-d10	19.04	212	13737m	0.21	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.44	128	2131	0.060	ng/ul#	90
5) 2-Methylnaphthalene	12.07	142	1122	0.046	ng/ul	98
7) Acenaphthylene	13.98	152	7521	0.193	ng/ul	98
8) Acenaphthene	14.32	153	1828	0.060	ng/ul	97
9) Fluorene	15.31	166	2286	0.066	ng/ul#	95
12) Phenanthrene	17.05	178	54115m	0.983	ng/ul	
13) Anthracene	17.14	178	11992m	0.247	ng/ul	
15) Fluoranthene	19.07	202	160690m	2.209	ng/ul	
17) Pyrene	19.44	202	153816	1.946	ng/ul	99
18) Benzo(a)anthracene	21.19	228	93120	1.290	ng/ul	97
19) Chrysene	21.24	228	103255	1.166	ng/ul	98
21) Benzo(b)fluoranthene	22.75	252	117785m	1.931	ng/ul	
22) Benzo(k)fluoranthene	22.78	252	37667m	0.545	ng/ul	
23) Benzo(a)pyrene	23.30	252	79594	1.240	ng/ul#	90
24) Indeno(1,2,3-cd)pyrene	25.57	276	49366	0.653	ng/ul#	94
25) Dibenzo(a,h)anthracene	25.58	278	11162	0.187	ng/ul#	80
26) Benzo(a,h,i)perylene	26.24	276	48642	0.702	ng/ul	100

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

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