

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN101119\
 Data File : BN008111.D
 Acq On : 12 Oct 2019 14:10
 Operator : JU
 Sample : K5178-02
 Misc :
 ALS Vial : 44 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 BEPD1

Manual Integrations
 APPROVED

mohammad
 10/15/2019 7:58:39 AM

Quant Time: Oct 14 15:48:43 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	2968	0.40	ng/ul	0.00
2) Naphthalene-d8	10.39	136	14351	0.40	ng/ul	-0.01
6) Acenaphthene-d10	14.26	164	9211	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.01	188	21416m	0.40	ng/ul	0.00
16) Chrysene-d12	21.21	240	20736m	0.40	ng/ul	0.00
20) Perylene-d12	23.40	264	18673	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.00	152	4621	0.19	ng/ul	0.00
14) Fluoranthene-d10	19.04	212	15766m	0.21	ng/ul	0.00
Target Compounds					Ovalue	
3) Naphthalene	10.44	128	849	0.021	ng/ul#	74
5) 2-Methylnaphthalene	12.07	142	587	0.021	ng/ul	99
7) Acenaphthylene	13.98	152	3743	0.085	ng/ul#	93
8) Acenaphthene	14.32	153	2032	0.060	ng/ul	97
9) Fluorene	15.31	166	2771	0.072	ng/ul#	97
12) Phenanthrene	17.05	178	119153m	1.900	ng/ul	
13) Anthracene	17.14	178	13559m	0.246	ng/ul	
15) Fluoranthene	19.07	202	442565m	5.344	ng/ul	
17) Pyrene	19.44	202	350596	4.246	ng/ul	100
18) Benzo(a)anthracene	21.20	228	164407	2.181	ng/ul	99
19) Chrysene	21.24	228	205012	2.216	ng/ul	99
21) Benzo(b)fluoranthene	22.75	252	255446m	4.202	ng/ul	
22) Benzo(k)fluoranthene	22.79	252	93369m	1.356	ng/ul	
23) Benzo(a)pyrene	23.31	252	147521	2.306	ng/ul#	85
24) Indeno(1,2,3-cd)pyrene	25.58	276	106010	1.407	ng/ul#	94
25) Dibenzo(a,h)anthracene	25.58	278	20791	0.349	ng/ul#	91
26) Benzo(a,h,i)perylene	26.24	276	103145	1.493	ng/ul	99

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

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