

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN112119\
 Data File : BN008698.D
 Acq On : 22 Nov 2019 04:52
 Operator : JU
 Sample : K5854-17
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 C0AD3

Manual Integrations
 APPROVED

mohammad
 11/22/2019 1:24:33 PM

Quant Time: Nov 22 07:37:45 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN111519.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Nov 22 07:12:34 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.61	152	4428	0.40	ng/ul	0.00
2) Naphthalene-d8	10.37	136	16182	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.23	164	9677	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.97	188	19491m	0.40	ng/ul	0.00
16) Chrysene-d12	21.17	240	13779	0.40	ng/ul	0.00
20) Perylene-d12	23.34	264	14600	0.40	ng/ul	0.01
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.96	152	5972	0.23	ng/ul	0.00
14) Fluoranthene-d10	19.01	212	14270	0.24	ng/ul	0.00
Target Compounds					Ovalue	
3) Naphthalene	10.42	128	104673	2.146	ng/ul	98
5) 2-Methylnaphthalene	12.04	142	71748	2.106	ng/ul	100
7) Acenaphthylene	13.95	152	25953	0.485	ng/ul	96
8) Acenaphthene	14.30	153	15120	0.381	ng/ul	99
9) Fluorene	15.29	166	21509	0.461	ng/ul#	94
12) Phenanthrene	17.02	178	242298	3.686	ng/ul	98
13) Anthracene	17.11	178	77252	1.244	ng/ul	99
15) Fluoranthene	19.04	202	394315	4.883	ng/ul	92
17) Pyrene	19.40	202	387680	7.013	ng/ul	98
18) Benzo(a)anthracene	21.16	228	212312	3.831	ng/ul	96
19) Chrysene	21.21	228	219023m	4.024	ng/ul	
21) Benzo(b)fluoranthene	22.70	252	310969m	5.165	ng/ul	
22) Benzo(k)fluoranthene	22.74	252	105392m	1.765	ng/ul	
23) Benzo(a)pyrene	23.25	252	193281	3.420	ng/ul#	87
24) Indeno(1,2,3-cd)pyrene	25.48	276	132082	1.999	ng/ul#	95
25) Dibenzo(a,h)anthracene	25.49	278	35042	0.657	ng/ul#	57
26) Benzo(a,h,i)perylene	26.13	276	123660	2.237	ng/ul	96

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

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