

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN112119\
 Data File : BN008699.D
 Acq On : 22 Nov 2019 05:28
 Operator : JU
 Sample : K5854-01
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 C0AB6

Manual Integrations
 APPROVED

mohammad
 11/22/2019 1:24:12 PM

Quant Time: Nov 22 07:40:17 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	4495	0.40	ng/ul	0.00
2) Naphthalene-d8	10.37	136	16792	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.23	164	9747	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.97	188	19328	0.40	ng/ul	0.00
16) Chrysene-d12	21.17	240	15451	0.40	ng/ul	0.00
20) Perylene-d12	23.34	264	16112	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.96	152	4565	0.17	ng/ul	0.00
14) Fluoranthene-d10	19.01	212	10314	0.17	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.42	128	17324	0.342	ng/ul	100
5) 2-Methylnaphthalene	12.04	142	20684	0.585	ng/ul	100
7) Acenaphthylene	13.95	152	2824	0.052	ng/ul#	88
8) Acenaphthene	14.30	153	6485	0.162	ng/ul	99
9) Fluorene	15.29	166	9908	0.211	ng/ul#	97
12) Phenanthrene	17.02	178	145951	2.239	ng/ul	98
13) Anthracene	17.11	178	31250	0.507	ng/ul	99
15) Fluoranthene	19.04	202	265626	3.317	ng/ul	94
17) Pyrene	19.40	202	225943	3.645	ng/ul	96
18) Benzo(a)anthracene	21.16	228	122170	1.966	ng/ul	98
19) Chrysene	21.21	228	140928	2.309	ng/ul	97
21) Benzo(b)fluoranthene	22.70	252	174577m	2.628	ng/ul	
22) Benzo(k)fluoranthene	22.73	252	54249m	0.823	ng/ul	
23) Benzo(a)pyrene	23.25	252	117612	1.886	ng/ul	94
24) Indeno(1,2,3-cd)pyrene	25.47	276	88517	1.214	ng/ul#	96
25) Dibenzo(a,h)anthracene	25.48	278	20571	0.349	ng/ul#	76
26) Benzo(a,h,i)perylene	26.12	276	92772	1.521	ng/ul	98

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

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