

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN112119\
 Data File : BN008702.D
 Acq On : 22 Nov 2019 07:16
 Operator : JU
 Sample : K5854-07
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 C0AC5

Manual Integrations
 APPROVED

mohammad
 11/22/2019 1:24:16 PM

Quant Time: Nov 22 08:20:48 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	4363	0.40	ng/ul	0.00
2) Naphthalene-d8	10.37	136	15947	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.23	164	9045	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.97	188	18298	0.40	ng/ul	0.00
16) Chrysene-d12	21.17	240	14798	0.40	ng/ul	0.00
20) Perylene-d12	23.33	264	15379	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.96	152	3349	0.13	ng/ul	0.00
14) Fluoranthene-d10	19.00	212	8010	0.14	ng/ul	0.00
Target Compounds					Ovalue	
3) Naphthalene	10.42	128	22663	0.471	ng/ul	99
5) 2-Methylnaphthalene	12.03	142	10717	0.319	ng/ul	100
7) Acenaphthylene	13.95	152	5647	0.113	ng/ul	96
8) Acenaphthene	14.29	153	4155	0.112	ng/ul	99
9) Fluorene	15.29	166	5938	0.136	ng/ul#	96
12) Phenanthrene	17.02	178	57101	0.925	ng/ul	98
13) Anthracene	17.11	178	17965	0.308	ng/ul	99
15) Fluoranthene	19.04	202	95604	1.261	ng/ul	94
17) Pyrene	19.40	202	89541	1.508	ng/ul	96
18) Benzo(a)anthracene	21.16	228	47504	0.798	ng/ul	95
19) Chrysene	21.21	228	49432m	0.846	ng/ul	
21) Benzo(b)fluoranthene	22.70	252	65858m	1.038	ng/ul	
22) Benzo(k)fluoranthene	22.73	252	24586m	0.391	ng/ul	
23) Benzo(a)pyrene	23.24	252	40342	0.678	ng/ul#	82
24) Indeno(1,2,3-cd)pyrene	25.47	276	29468	0.423	ng/ul#	97
25) Dibenzo(a,h)anthracene	25.48	278	7892	0.140	ng/ul#	40
26) Benzo(a,h,i)perylene	26.12	276	27551	0.473	ng/ul#	91

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

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