

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN121119\
 Data File : BN008934.D
 Acq On : 12 Dec 2019 01:19
 Operator : JU
 Sample : K6088-19
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 C0BP7

Manual Integrations
 APPROVED

mohammad
 12/12/2019 11:32:46 AM

Quant Time: Dec 12 02:54:24 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 : Thu Dec 12 01:20:09 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.56	152	2652	0.40	ng/ul	0.00
2) Naphthalene-d8	10.31	136	11390	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.19	164	7280	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.93	188	16959	0.40	ng/ul	0.00
16) Chrysene-d12	21.13	240	14464	0.40	ng/ul	0.00
20) Perylene-d12	23.27	264	15446	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.91	152	3258	0.17	ng/ul	0.00
14) Fluoranthene-d10	18.96	212	10053	0.19	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.36	128	9228	0.269	ng/ul	99
5) 2-Methylnaphthalene	11.98	142	2939	0.123	ng/ul	100
7) Acenaphthylene	13.90	152	53288	1.324	ng/ul	100
8) Acenaphthene	14.25	153	5570	0.186	ng/ul	98
9) Fluorene	15.24	166	13872	0.395	ng/ul	99
12) Phenanthrene	16.97	178	347442	6.075	ng/ul	99
13) Anthracene	17.06	178	66731	1.235	ng/ul	100
15) Fluoranthene	19.00	202	1133332	16.129	ng/ul	98
17) Pyrene	19.36	202	968934	16.698	ng/ul	98
18) Benzo(a)anthracene	21.11	228	493728	8.488	ng/ul	98
19) Chrysene	21.17	228	476496	8.340	ng/ul	98
21) Benzo(b)fluoranthene	22.64	252	675356m	10.603	ng/ul	
22) Benzo(k)fluoranthene	22.68	252	270650m	4.285	ng/ul	
23) Benzo(a)pyrene	23.19	252	500765m	8.376	ng/ul	
24) Indeno(1,2,3-cd)pyrene	25.38	276	313819	4.488	ng/ul	90
25) Dibenzo(a,h)anthracene	25.39	278	72084	1.277	ng/ul	99
26) Benzo(a,h,i)perylene	26.02	276	298854	5.110	ng/ul	96

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

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