

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN121119\
 Data File : BN008937.D
 Acq On : 12 Dec 2019 10:08
 Operator : JU
 Sample : K6088-19DL 5X
 Misc :
 ALS Vial : 28 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 C0BP7DL

Manual Integrations
 APPROVED

mohammad
 12/13/2019 8:07:36 AM

Quant Time: Dec 12 11:17:19 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 00:49:42 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.55	152	2757	0.40	ng/ul	-0.01
2) Naphthalene-d8	10.30	136	11381	0.40	ng/ul	-0.01
6) Acenaphthene-d10	14.18	164	7404	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.92	188	17110	0.40	ng/ul	0.00
16) Chrysene-d12	21.13	240	14643	0.40	ng/ul	0.00
20) Perylene-d12	23.28	264	14405	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.90	152	710	0.04	ng/ul	-0.01
14) Fluoranthene-d10	18.96	212	2130	0.04	ng/ul	0.00
Target Compounds					Ovalue	
3) Naphthalene	10.35	128	1954	0.057	ng/ul#	91
5) 2-Methylnaphthalene	11.97	142	652	0.027	ng/ul	97
7) Acenaphthylene	13.89	152	11228	0.274	ng/ul	99
8) Acenaphthene	14.24	153	1258	0.041	ng/ul	97
9) Fluorene	15.23	166	2904	0.081	ng/ul#	97
12) Phenanthrene	16.97	178	72099	1.249	ng/ul	99
13) Anthracene	17.06	178	14097	0.259	ng/ul	99
15) Fluoranthene	18.99	202	241559	3.407	ng/ul	100
17) Pyrene	19.35	202	210003	3.575	ng/ul	99
18) Benzo(a)anthracene	21.12	228	98290	1.669	ng/ul	98
19) Chrysene	21.17	228	97382	1.684	ng/ul	98
21) Benzo(b)fluoranthene	22.64	252	137864m	2.321	ng/ul	
22) Benzo(k)fluoranthene	22.68	252	57773m	0.981	ng/ul	
23) Benzo(a)pyrene	23.19	252	100212m	1.797	ng/ul	
24) Indeno(1,2,3-cd)pyrene	25.38	276	66187	1.015	ng/ul	91
25) Dibenzo(a,h)anthracene	25.39	278	15012	0.285	ng/ul	96
26) Benzo(a,h,i)perylene	26.02	276	63937	1.172	ng/ul	97

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

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