

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN121119\
 Data File : BN008920.D
 Acq On : 11 Dec 2019 16:59
 Operator : JU
 Sample : K6088-05
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 C0BN2

Manual Integrations
 APPROVED

mohammad
 12/12/2019 11:27:11 AM

Quant Time: Dec 12 01:23:09 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 : Wed Dec 11 10:35:08 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.56	152	2058	0.40	ng/ul	0.00
2) Naphthalene-d8	10.31	136	8970	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.19	164	5881	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.93	188	13481	0.40	ng/ul	0.00
16) Chrysene-d12	21.13	240	12434	0.40	ng/ul	0.00
20) Perylene-d12	23.28	264	13863	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.91	152	2654	0.18	ng/ul	0.00
14) Fluoranthene-d10	18.96	212	8599	0.21	ng/ul	0.00
Target Compounds					Ovalue	
3) Naphthalene	10.36	128	18245	0.675	ng/ul	100
5) 2-Methylnaphthalene	11.98	142	14751	0.781	ng/ul	100
7) Acenaphthylene	13.90	152	12994	0.400	ng/ul	98
8) Acenaphthene	14.25	153	931	0.039	ng/ul	95
9) Fluorene	15.24	166	2095	0.074	ng/ul#	86
12) Phenanthrene	16.97	178	50218	1.105	ng/ul	99
13) Anthracene	17.06	178	14504	0.338	ng/ul	98
15) Fluoranthene	19.00	202	98172	1.758	ng/ul	99
17) Pyrene	19.36	202	94980	1.904	ng/ul	99
18) Benzo(a)anthracene	21.12	228	84967	1.699	ng/ul	96
19) Chrysene	21.17	228	95909	1.953	ng/ul	98
21) Benzo(b)fluoranthene	22.65	252	137117m	2.399	ng/ul	
22) Benzo(k)fluoranthene	22.69	252	50908m	0.898	ng/ul	
23) Benzo(a)pyrene	23.19	252	85841	1.600	ng/ul#	94
24) Indeno(1,2,3-cd)pyrene	25.39	276	52882	0.843	ng/ul#	91
25) Dibenzo(a,h)anthracene	25.39	278	14239	0.281	ng/ul	93
26) Benzo(a,h,i)perylene	26.03	276	45011	0.857	ng/ul	97

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

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