

Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN121119\
 Data File : BN008944.D
 Acq On : 12 Dec 2019 14:17
 Operator : JU
 Sample : K6089-01
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 C0BP8

Manual Integrations
 APPROVED

mohammad
 12/13/2019 9:07:31 AM

Quant Time: Dec 12 15:13:40 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	7.56	152	2213	0.40	ng/ul	0.00
2) Naphthalene-d8	10.31	136	9547	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.18	164	6116	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.93	188	14209m	0.40	ng/ul	0.00
16) Chrysene-d12	21.13	240	12617	0.40	ng/ul	0.00
20) Perylene-d12	23.28	264	13343	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.91	152	3344	0.21	ng/ul	0.00
14) Fluoranthene-d10	18.96	212	9825	0.23	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	10.36	128	588	0.020	ng/ul#	72
7) Acenaphthylene	13.90	152	1926	0.057	ng/ul#	91
12) Phenanthrene	16.97	178	7135	0.149	ng/ul	98
13) Anthracene	17.06	178	1625	0.036	ng/ul#	89
15) Fluoranthene	18.99	202	25497	0.433	ng/ul	98
17) Pyrene	19.36	202	24620	0.486	ng/ul	98
18) Benzo(a)anthracene	21.12	228	11254	0.222	ng/ul	97
19) Chrysene	21.17	228	11946	0.240	ng/ul	98
21) Benzo(b)fluoranthene	22.64	252	17494m	0.318	ng/ul	
22) Benzo(k)fluoranthene	22.68	252	7405m	0.136	ng/ul	
23) Benzo(a)pyrene	23.19	252	12384	0.240	ng/ul	94
24) Indeno(1,2,3-cd)pyrene	25.38	276	8429	0.140	ng/ul#	94
25) Dibenzo(a,h)anthracene	25.38	278	1863	0.038	ng/ul#	63
26) Benzo(g,h,i)perylene	26.02	276	8264	0.164	ng/ul	98

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

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