

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN121319\
 Data File : BN008982.D
 Acq On : 14 Dec 2019 00:22
 Operator : JU
 Sample : K4649-02MS
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 A5167MS

Quant Time: Dec 14 02:36:34 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 Dec 13 23:22:26 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.56	152	2381	0.40	ng/ul	0.00
2) Naphthalene-d8	10.31	136	10550	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.18	164	7064	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.92	188	15862	0.40	ng/ul	0.00
16) Chrysene-d12	21.12	240	12712	0.40	ng/ul	0.00
20) Perylene-d12	23.27	264	10638	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.90	152	7372	0.43	ng/ul	0.00
14) Fluoranthene-d10	18.96	212	18811	0.39	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.36	128	12737	0.400	ng/ul	99
5) 2-Methylnaphthalene	11.98	142	9591	0.432	ng/ul	100
7) Acenaphthylene	13.89	152	14224	0.364	ng/ul	100
8) Acenaphthene	14.24	153	10926	0.377	ng/ul	98
9) Fluorene	15.24	166	12866	0.378	ng/ul	99
11) Pentachlorophenol	16.59	266	1642	0.274	ng/ul	99
12) Phenanthrene	16.97	178	21753	0.407	ng/ul	100
13) Anthracene	17.06	178	19920	0.394	ng/ul	99
15) Fluoranthene	18.99	202	25211	0.384	ng/ul	99
17) Pyrene	19.35	202	25303	0.496	ng/ul	98
18) Benzo(a)anthracene	21.11	228	23386	0.457	ng/ul	100
19) Chrysene	21.16	228	22835	0.455	ng/ul	100
21) Benzo(b)fluoranthene	22.64	252	19712	0.449	ng/ul	99
22) Benzo(k)fluoranthene	22.68	252	19654	0.452	ng/ul	99
23) Benzo(a)pyrene	23.18	252	17271	0.419	ng/ul	99
24) Indeno(1,2,3-cd)pyrene	25.37	276	17598	0.365	ng/ul#	99
25) Dibenzo(a,h)anthracene	25.39	278	14307	0.368	ng/ul	99
26) Benzo(g,h,i)perylene	26.01	276	14621	0.363	ng/ul	98

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

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