

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN121319\
 Data File : BN008983.D
 Acq On : 14 Dec 2019 00:58
 Operator : JU
 Sample : K4649-03MSD
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 A5167MSD

Quant Time: Dec 14 02:36:48 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	7.55	152	2720	0.40	ng/ul	0.00
2) Naphthalene-d8	10.31	136	11434	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.18	164	7575	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.92	188	16925	0.40	ng/ul	0.00
16) Chrysene-d12	21.13	240	13535	0.40	ng/ul	0.00
20) Perylene-d12	23.27	264	11311	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.90	152	8027	0.43	ng/ul	0.00
14) Fluoranthene-d10	18.96	212	20120	0.39	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.36	128	13773	0.400	ng/ul	100
5) 2-Methylnaphthalene	11.98	142	10322	0.429	ng/ul	100
7) Acenaphthylene	13.89	152	15080	0.360	ng/ul	99
8) Acenaphthene	14.24	153	11675	0.376	ng/ul	99
9) Fluorene	15.24	166	13772	0.377	ng/ul	100
11) Pentachlorophenol	16.59	266	1735	0.271	ng/ul	98
12) Phenanthrene	16.97	178	23148	0.406	ng/ul	99
13) Anthracene	17.06	178	21066	0.391	ng/ul	99
15) Fluoranthene	18.99	202	26718	0.381	ng/ul	99
17) Pyrene	19.35	202	26723	0.492	ng/ul	99
18) Benzo(a)anthracene	21.11	228	24396	0.448	ng/ul	100
19) Chrysene	21.16	228	24276	0.454	ng/ul	100
21) Benzo(b)fluoranthene	22.64	252	21049	0.451	ng/ul	99
22) Benzo(k)fluoranthene	22.68	252	20754	0.449	ng/ul	100
23) Benzo(a)pyrene	23.18	252	18447	0.421	ng/ul	99
24) Indeno(1,2,3-cd)pyrene	25.38	276	19201	0.375	ng/ul	99
25) Dibenzo(a,h)anthracene	25.39	278	15595	0.377	ng/ul	98
26) Benzo(g,h,i)perylene	26.02	276	15974	0.373	ng/ul	98

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

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