

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN121718\
 Data File : BN004067.D
 Acq On : 18 Dec 2018 09:39
 Operator : JU/SJ
 Sample : J6270-15
 Misc :
 ALS Vial : 34 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 F9E14

Quant Time: Dec 18 11:34:45 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-SIM-BN111918.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Dec 18 06:35:27 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.84	152	1651	0.40	ng/ul	0.00
2) Naphthalene-d8	10.63	136	8435	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.47	164	5293	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.31	188	687184	0.40	ng/ul	0.10
16) Chrysene-d12	0.00	240	0	0.00	ng/ul	-21.40
20) Perylene-d12	23.73	264	15309	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.22	152	2868	0.22	ng/ul	0.00
14) Fluoranthene-d10	0.00	212	0	0.00	ng/ul	
Target Compounds				Qvalue		
7) Acenaphthylene	14.19	152	1613	0.068	ng/ul#	89
21) Benzo(b)fluoranthene	23.02	252	13991	0.221	ng/ul#	79
22) Benzo(k)fluoranthene	23.02	252	13979	0.227	ng/ul#	79
23) Benzo(a)pyrene	23.62	252	4459	0.081	ng/ul#	63
24) Indeno(1,2,3-cd)pyrene	26.09	276	4530	0.072	ng/ul#	100
25) Dibenzo(a,h)anthracene	26.08	278	1299	0.026	ng/ul#	1
26) Benzo(g,h,i)perylene	26.82	276	4944	0.090	ng/ul#	73

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

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Quant Time: Dec 18 11:34:45 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-SIM-BN11918.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Dec 18 06:35:27 2018
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