

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN061119\
 Data File : BN006264.D
 Acq On : 12 Jun 2019 02:08
 Operator : JU/SJ
 Sample : K3215-08
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 DB8J2

Manual Integrations
 APPROVED

mohammad
 6/12/2019 8:19:28 AM

Quant Time: Jun 12 03:56:55 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN053019.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sun Jun 09 05:10:30 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.64	152	6097	0.40	ng/ul	0.00
2) Naphthalene-d8	10.42	136	24074	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.30	164	14477	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.05	188	25059m	0.40	ng/ul	0.00
16) Chrysene-d12	21.26	240	15075m	0.40	ng/ul	0.00
20) Perylene-d12	23.49	264	26393	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.02	152	11606	0.31	ng/ul	0.00
14) Fluoranthene-d10	19.09	212	20237m	0.30	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.47	128	132279	1.991	ng/ul	100
5) 2-Methylnaphthalene	12.09	142	194516	4.202	ng/ul	100
7) Acenaphthylene	14.01	152	3920	0.049	ng/ul#	1
8) Acenaphthene	14.36	153	13142	0.235	ng/ul#	81
9) Fluorene	15.35	166	16735	0.256	ng/ul	89
11) Pentachlorophenol	16.71	266	505393m	79.371	ng/ul	
12) Phenanthrene	17.09	178	19734	0.245	ng/ul	96

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

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN061119\
Data File : BN006264.D
Acq On : 12 Jun 2019 02:08
Operator : JU/SJ
Sample : K3215-08
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
DB8J2

Manual Integrations
APPROVED

mohammad
6/12/2019 8:19:28 AM

Quant Time: Jun 12 03:56:55 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN053019.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Sun Jun 09 05:10:30 2019
Response via : Initial Calibration

