

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN120418\
 Data File : BN003701.D
 Acq On : 04 Dec 2018 15:46
 Operator : JU/SJ
 Sample : J6137-07DL 2X
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 F9J00DL

Manual Integrations
 APPROVED

mohammad
 12/6/2018 10:34:07 AM

Quant Time: Dec 04 16:47:14 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 04 11:17:33 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.84	152	5124	0.40	ng/ul	0.00
2) Naphthalene-d8	10.63	136	20870	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.47	164	11518	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.22	188	25585m	0.40	ng/ul	0.00
16) Chrysene-d12	21.40	240	22514m	0.40	ng/ul	0.00
20) Perylene-d12	23.71	264	23671	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.22	152	5117	0.16	ng/ul	0.00
14) Fluoranthene-d10	19.24	212	13110m	0.18	ng/ul	0.00
Target Compounds						
3) Naphthalene	10.68	128	1154140	19.688	ng/ul	100
5) 2-Methylnaphthalene	12.29	142	36344	0.888	ng/ul	100
7) Acenaphthylene	14.20	152	2910	0.056	ng/ul#	94
8) Acenaphthene	14.54	153	12170	0.268	ng/ul	99
9) Fluorene	15.52	166	7113	0.134	ng/ul	98
11) Pentachlorophenol	16.86	266	320m	0.069	ng/ul	
12) Phenanthrene	17.26	178	4548	0.054	ng/ul	96

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

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN120418\
Data File : BN003701.D
Acq On : 04 Dec 2018 15:46
Operator : JU/SJ
Sample : J6137-07DL 2X
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
F9J00DL

Manual Integrations
APPROVED

mohammad
12/6/2018 10:34:07 AM

Quant Time: Dec 04 16:47:14 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 04 11:17:33 2018
Response via : Initial Calibration

