

Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN120518\
 Data File : BN003721.D
 Acq On : 05 Dec 2018 21:20
 Operator : JU/SJ
 Sample : J6170-07
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 F9J04

Manual Integrations
 APPROVED

mohammad
 12/6/2018 3:58:38 PM

Quant Time: Dec 06 03:21:31 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 : Thu Dec 06 02:51:43 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.84	152	4887	0.40	ng/ul	0.00
2) Naphthalene-d8	10.63	136	20126	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.47	164	12298	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.21	188	24927m	0.40	ng/ul	0.00
16) Chrysene-d12	21.39	240	20921m	0.40	ng/ul	0.00
20) Perylene-d12	23.70	264	15721m	0.40	ng/ul	-0.01
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.22	152	11432	0.37	ng/ul	0.00
14) Fluoranthene-d10	19.24	212	27080m	0.37	ng/ul	0.00
Target Compounds						
3) Naphthalene	10.69	128	6658918	117.790	ng/ul	98
5) 2-Methylnaphthalene	12.29	142	243967	6.180	ng/ul	100
7) Acenaphthylene	14.20	152	21334	0.385	ng/ul	98
8) Acenaphthene	14.53	153	112713	2.323	ng/ul	99
9) Fluorene	15.52	166	61665	1.090	ng/ul	99
11) Pentachlorophenol	16.86	266	9837m	2.191	ng/ul	
12) Phenanthrene	17.26	178	37927	0.464	ng/ul	99
13) Anthracene	17.35	178	2815m	0.042	ng/ul	

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

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-SIM-BN11918.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Dec 06 02:51:43 2018
Response via : Initial Calibration

