

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN120418\
 Data File : BN003697.D
 Acq On : 04 Dec 2018 12:40
 Operator : JU/SJ
 Sample : J6137-07
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 F9J00

Manual Integrations
 APPROVED

mohammad
 12/6/2018 10:33:56 AM

Quant Time: Dec 04 15:12:42 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	5595	0.40	ng/ul	0.00
2) Naphthalene-d8	10.63	136	22958	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.47	164	13193	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.22	188	28905m	0.40	ng/ul	0.00
16) Chrysene-d12	21.40	240	25917m	0.40	ng/ul	0.00
20) Perylene-d12	23.70	264	20980m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.22	152	10539	0.30	ng/ul	0.00
14) Fluoranthene-d10	19.24	212	26345m	0.31	ng/ul	0.00
Target Compounds						
3) Naphthalene	10.68	128	2263941	35.107	ng/ul	100
5) 2-Methylnaphthalene	12.29	142	75144	1.669	ng/ul	100
7) Acenaphthylene	14.19	152	5899	0.099	ng/ul	97
8) Acenaphthene	14.53	153	25045	0.481	ng/ul	100
9) Fluorene	15.52	166	14618	0.241	ng/ul	98
11) Pentachlorophenol	16.86	266	685m	0.132	ng/ul	
12) Phenanthrene	17.26	178	9302	0.098	ng/ul	99

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

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