

Data Path : U:\HPCHEM1\BNA E\DATA\BE020218\
 Data File : BE095398.D
 Acq On : 2 Feb 2018 15:49
 Operator : SJ/JU
 Sample : J1245-06DL 5X
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 F6B67DL

Quant Time: Feb 05 15:52:51 2018
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\SOM-EPA-SIM-BE012618.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Feb 05 04:46:10 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.44	152	1905	0.40	ng/ul	0.00
2) Naphthalene-d8	11.26	136	5791	0.40	ng/ul	0.00
6) Acenaphthene-d10	15.02	164	3818	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.71	188	6989	0.40	ng/ul	0.00
16) Chrysene-d12	21.80	240	7210	0.40	ng/ul	0.00
20) Perylene-d12	24.44	264	7368	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.81	152	766	0.08	ng/ul	0.00
14) Fluoranthene-d10	19.69	212	1436	0.05	ng/ul	0.00
Target Compounds				Ovalue		
5) 2-Methylnaphthalene	12.88	142	36690	3.388	ng/ul	100
8) Acenaphthene	15.08	153	3165	0.222	ng/ul#	61
9) Fluorene	16.03	166	5520	0.376	ng/ul#	67
11) Pentachlorophenol	17.36	266	2064	1.525	ng/ul	99
12) Phenanthrene	17.75	178	23331	0.993	ng/ul#	94
15) Fluoranthene	19.72	202	2900	0.095	ng/ul	83
17) Pyrene	20.08	202	5910	0.244	ng/ul#	86

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

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