

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE040918\
 Data File : BE095964.D
 Acq On : 9 Apr 2018 13:14
 Operator : SJ/JU
 Sample : J2135-08
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 X2C46

Quant Time: Apr 10 03:57:21 2018
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SOM-EPA-SIM-BE040618.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Apr 10 03:47:21 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.39	152	1980	0.40	ng/ul	0.00
2) Naphthalene-d8	11.21	136	5255	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.97	164	2918	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.68	188	7053	0.40	ng/ul	0.00
16) Chrysene-d12	21.77	240	8946	0.40	ng/ul	0.00
20) Perylene-d12	24.39	264	7188	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.77	152	2588	0.26	ng/ul	0.00
14) Fluoranthene-d10	19.66	212	9493	0.33	ng/ul	0.00
Target Compounds					Qvalue	
3) Naphthalene	11.26	128	6929	0.472	ng/ul#	88
5) 2-Methylnaphthalene	12.84	142	5200	0.523	ng/ul	99
7) Acenaphthylene	14.70	152	8521	0.563	ng/ul	97
8) Acenaphthene	15.04	153	6995	0.645	ng/ul	95
11) Pentachlorophenol	17.34	266	2171	1.454	ng/ul	96
13) Anthracene	17.81	178	10889	0.526	ng/ul#	91
17) Pyrene	20.05	202	26203	0.862	ng/ul#	79
19) Chrysene	21.82	228	16778	0.645	ng/ul	97
21) Benzo(b)fluoranthene	23.58	252	19283	0.829	ng/ul	85
22) Benzo(k)fluoranthene	23.64	252	28289	1.221	ng/ul#	85
25) Dibenzo(a,h)anthracene	27.21	278	16965	0.845	ng/ul#	88

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

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