

Data Path : Z:\SVOASRV\HPCHEM1\BNA_E\DATA\BE082918\
 Data File : BE097393.D
 Acq On : 29 Aug 2018 17:24
 Operator : SJ/JU
 Sample : J4596-07MS
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 A3Z29MS

Manual Integrations
 APPROVED

Sohil
 8/30/2018 5:36:45 PM

Quant Time: Aug 29 18:18:52 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_E\METHODS\SOM-EPA-SIM-BE082918.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Aug 29 13:18:39 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.38	152	841	0.40	ng/ul	0.00
2) Naphthalene-d8	10.12	136	4212	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.01	164	2582	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.76	188	7570	0.40	ng/ul	0.00
16) Chrysene-d12	20.97	240	9885	0.40	ng/ul	0.00
20) Perylene-d12	23.20	264	9793	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.72	152	883	0.13	ng/ul	0.00
14) Fluoranthene-d10	18.80	212	3876	0.16	ng/ul	0.00
Target Compounds					Qvalue	
8) Acenaphthene	14.07	153	931	0.115	ng/ul#	94
11) Pentachlorophenol	16.44	266	19	0.025	ng/ul#	70
12) Phenanthrene	16.80	178	716	0.037	ng/ul#	83
15) Fluoranthene	18.83	202	1152	0.046	ng/ul	86
17) Pyrene	19.20	202	4925	0.185	ng/ul#	81
18) Benzo(a)anthracene	20.95	228	684	0.027	ng/ul#	65
19) Chrysene	20.98	228	1512	0.052	ng/ul#	71
21) Benzo(b)fluoranthene	22.53	252	1512m	0.058	ng/ul	
22) Benzo(k)fluoranthene	22.57	252	752m	0.026	ng/ul	
23) Benzo(a)pyrene	23.10	252	529	0.022	ng/ul#	16
26) Benzo(g,h,i)perylene	26.20	276	651	0.024	ng/ul#	66

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

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