

Data Path : Z:\SVOASRV\HPCHEM1\BNA_E\DATA\BE082918\
 Data File : BE097389.D
 Acq On : 29 Aug 2018 14:52
 Operator : SJ/JU
 Sample : J4596-03
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 A3YT5

Quant Time: Aug 29 15:36:06 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	717	0.40	ng/ul	0.00
2) Naphthalene-d8	10.12	136	3599	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.01	164	2124	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.76	188	6605	0.40	ng/ul	0.00
16) Chrysene-d12	20.97	240	9200	0.40	ng/ul	0.00
20) Perylene-d12	23.20	264	9043	0.40	ng/ul	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
4) 2-Methylnaphthalene-d10	11.72	152	2032	0.36	ng/ul	0.00
14) Fluoranthene-d10	18.80	212	10066	0.46	ng/ul	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
3) Naphthalene	10.17	128	2588	0.295	ng/ul#	96
5) 2-Methylnaphthalene	11.79	142	1573	0.284	ng/ul	98
7) Acenaphthylene	13.73	152	2346	0.303	ng/ul	99
9) Fluorene	15.07	166	3093	0.406	ng/ul	95
11) Pentachlorophenol	16.44	266	76	0.117	ng/ul	87
12) Phenanthrene	16.80	178	12188	0.717	ng/ul#	89
13) Anthracene	16.89	178	1857	0.130	ng/ul	94
15) Fluoranthene	18.83	202	11258	0.515	ng/ul	84
17) Pyrene	19.20	202	13126	0.530	ng/ul#	80
18) Benzo(a)anthracene	20.95	228	16243	0.687	ng/ul#	85
19) Chrysene	21.01	228	11333	0.416	ng/ul	98
21) Benzo(b)fluoranthene	22.53	252	7682	0.317	ng/ul	85
22) Benzo(k)fluoranthene	22.57	252	13444	0.498	ng/ul#	73
24) Indeno(1,2,3-cd)pyrene	25.49	276	11075	0.381	ng/ul#	80
25) Dibenzo(a,h)anthracene	25.51	278	10359	0.421	ng/ul#	92
26) Benzo(g,h,i)perylene	26.19	276	14521	0.587	ng/ul#	91

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

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