

Data Path : Z:\SVOASRV\HPCHEM1\BNA_E\DATA\BE083118\
 Data File : BE097403.D
 Acq On : 31 Aug 2018 13:19
 Operator : SJ/JU
 Sample : J4596-07MS
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 A3Z29MS

Quant Time: Aug 31 14:29:02 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 : Fri Aug 31 11:16:59 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.38	152	741	0.40	ng/ul	0.00
2) Naphthalene-d8	10.12	136	3815	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.01	164	2374	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.76	188	7263	0.40	ng/ul	0.00
16) Chrysene-d12	20.97	240	9365	0.40	ng/ul	0.00
20) Perylene-d12	23.20	264	9398	0.40	ng/ul	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
4) 2-Methylnaphthalene-d10	11.72	152	1023	0.17	ng/ul	0.00
14) Fluoranthene-d10	18.80	212	4724	0.20	ng/ul	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
3) Naphthalene	10.17	128	231	0.025	ng/ul#	56
8) Acenaphthene	14.07	153	1175	0.158	ng/ul#	95
11) Pentachlorophenol	16.43	266	296	0.414	ng/ul	97
12) Phenanthrene	16.80	178	1352	0.072	ng/ul#	92
13) Anthracene	16.90	178	529	0.034	ng/ul#	79
15) Fluoranthene	18.83	202	2321	0.097	ng/ul	84
17) Pyrene	19.20	202	7069	0.281	ng/ul#	80
18) Benzo(a)anthracene	20.95	228	1485	0.062	ng/ul#	76
19) Chrysene	21.00	228	3072	0.111	ng/ul#	80
21) Benzo(b)fluoranthene	22.53	252	3154	0.125	ng/ul#	69
22) Benzo(k)fluoranthene	22.57	252	1538	0.055	ng/ul#	1
23) Benzo(a)pyrene	23.10	252	1224	0.053	ng/ul#	51
24) Indeno(1,2,3-cd)pyrene	25.49	276	1475	0.049	ng/ul#	89
25) Dibenzo(a,h)anthracene	25.51	278	820	0.032	ng/ul#	1
26) Benzo(g,h,i)perylene	26.19	276	1497	0.058	ng/ul#	81

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

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