

Data Path : Z:\svoasrv\HPCHEM1\BNA_E\Data\BE083118\
 Data File : BE097404.D
 Acq On : 31 Aug 2018 13:58
 Operator : SJ/JU
 Sample : J4596-08MSD
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 A3Z29MSD

Quant Time: Aug 31 14:48:35 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	728	0.40	ng/ul	0.00
2) Naphthalene-d8	10.12	136	3729	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.01	164	2427	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.76	188	7405	0.40	ng/ul	0.00
16) Chrysene-d12	20.97	240	9579	0.40	ng/ul	0.00
20) Perylene-d12	23.20	264	9648	0.40	ng/ul	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
4) 2-Methylnaphthalene-d10	11.72	152	1018	0.18	ng/ul	0.00
14) Fluoranthene-d10	18.80	212	4856	0.20	ng/ul	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
3) Naphthalene	10.17	128	233	0.026	ng/ul#	55
8) Acenaphthene	14.07	153	1217	0.160	ng/ul	95
11) Pentachlorophenol	16.43	266	307	0.421	ng/ul	98
12) Phenanthrene	16.80	178	1402	0.074	ng/ul#	92
13) Anthracene	16.89	178	542	0.034	ng/ul#	81
15) Fluoranthene	18.83	202	2342	0.096	ng/ul	85
17) Pyrene	19.20	202	7208	0.280	ng/ul#	81
18) Benzo(a)anthracene	20.95	228	1550	0.063	ng/ul#	74
19) Chrysene	20.98	228	2899	0.102	ng/ul#	76
21) Benzo(b)fluoranthene	22.53	252	3244	0.126	ng/ul#	73
22) Benzo(k)fluoranthene	22.57	252	1493	0.052	ng/ul#	1
23) Benzo(a)pyrene	23.10	252	1195	0.050	ng/ul#	52
24) Indeno(1,2,3-cd)pyrene	25.48	276	1556	0.050	ng/ul#	86
25) Dibenzo(a,h)anthracene	25.51	278	823	0.031	ng/ul#	1
26) Benzo(g,h,i)perylene	26.19	276	1606	0.061	ng/ul#	82

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

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