

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE080118\
 Data File : BE097126.D
 Acq On : 1 Aug 2018 12:27
 Operator : SJ/JU
 Sample : SST00.173
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SST00.173

Quant Time: Aug 01 14:23:35 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\SOM-EPA-SIM-BE080118.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Aug 01 14:22:12 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.39	152	1485	0.40	ng/ul	0.00
2) Naphthalene-d8	10.14	136	6770	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.02	164	3765	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.77	188	10415	0.40	ng/ul	0.00
16) Chrysene-d12	20.98	240	14422	0.40	ng/ul	0.00
20) Perylene-d12	23.21	264	15307	0.40	ng/ul	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
4) 2-Methylnaphthalene-d10	11.73	152	1065	0.10	ng/ul	0.00
14) Fluoranthene-d10	18.81	212	3744	0.14	ng/ul	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Ovalue
3) Naphthalene	10.18	128	1683	0.103	ng/ul#	93
5) 2-Methylnaphthalene	11.81	142	1075	0.096	ng/ul	100
7) Acenaphthylene	13.73	152	1633	0.106	ng/ul	96
8) Acenaphthene	14.08	153	1171	0.092	ng/ul#	90
9) Fluorene	15.08	166	1371	0.104	ng/ul	97
12) Phenanthrene	16.81	178	2709	0.101	ng/ul#	91
13) Anthracene	16.90	178	2447	0.107	ng/ul	93
15) Fluoranthene	18.84	202	3682	0.129	ng/ul	85
17) Pyrene	19.21	202	3913	0.083	ng/ul#	87
18) Benzo(a)anthracene	20.96	228	4210	0.124	ng/ul#	85
19) Chrysene	21.01	228	4279	0.097	ng/ul	96
21) Benzo(b)fluoranthene	22.53	252	4057	0.097	ng/ul#	25
22) Benzo(k)fluoranthene	22.58	252	4572	0.100	ng/ul#	9
23) Benzo(a)pyrene	23.11	252	4208	0.119	ng/ul#	2
24) Indeno(1,2,3-cd)pyrene	25.50	276	4899	0.134	ng/ul#	82
25) Dibenzo(a,h)anthracene	25.53	278	4040	0.139	ng/ul#	49
26) Benzo(a,h,i)perylene	26.21	276	4182	0.117	ng/ul#	79

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

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