

Data Path : Z:\SVOASRV\HPCHEM1\BNA_E\DATA\BE082918\
 Data File : BE097384.D
 Acq On : 29 Aug 2018 11:17
 Operator : SJ/JU
 Sample : SST00.816
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SST00.816

Quant Time: Aug 29 13:14:16 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 13:07:59 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.38	152	830	0.40	ng/ul	0.00
2) Naphthalene-d8	10.12	136	3967	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.01	164	2223	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.76	188	6284	0.40	ng/ul	0.00
16) Chrysene-d12	20.97	240	8085	0.40	ng/ul	0.00
20) Perylene-d12	23.20	264	7258	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.72	152	4930	0.78	ng/ul	0.00
14) Fluoranthene-d10	18.80	212	16626	0.99	ng/ul	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
3) Naphthalene	10.17	128	7582	0.783	ng/ul#	92
5) 2-Methylnaphthalene	11.79	142	4971	0.796	ng/ul	97
7) Acenaphthylene	13.73	152	6543	0.652	ng/ul	96
8) Acenaphthene	14.07	153	5506	0.762	ng/ul	96
9) Fluorene	15.07	166	6506	0.828	ng/ul#	95
11) Pentachlorophenol	16.43	266	461	1.130	ng/ul	98
12) Phenanthrene	16.80	178	12840	0.791	ng/ul#	90
13) Anthracene	16.89	178	11141	0.819	ng/ul#	91
15) Fluoranthene	18.83	202	16792	0.973	ng/ul	84
17) Pyrene	19.20	202	16195	0.663	ng/ul#	79
18) Benzo(a)anthracene	20.95	228	16425	0.819	ng/ul#	85
19) Chrysene	21.01	228	18730	0.783	ng/ul	97
21) Benzo(b)fluoranthene	22.53	252	16009	0.815	ng/ul	82
22) Benzo(k)fluoranthene	22.57	252	17333	0.710	ng/ul#	85
23) Benzo(a)pyrene	23.10	252	14658	0.807	ng/ul#	76
24) Indeno(1,2,3-cd)pyrene	25.49	276	18597	0.824	ng/ul#	76
25) Dibenzo(a,h)anthracene	25.51	278	15844	0.868	ng/ul#	86
26) Benzo(g,h,i)perylene	26.19	276	16026	0.886	ng/ul#	91

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

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