

Data Path : Z:\SVOASRV\HPCHEM1\BNA_E\DATA\BE082918\
 Data File : BE097387.D
 Acq On : 29 Aug 2018 13:37
 Operator : SJ/JU
 Sample : SSTDICV0.4
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SICV19

Quant Time: Aug 29 14:21:11 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	849	0.40	ng/ul	0.00
2) Naphthalene-d8	10.12	136	4098	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.01	164	2264	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.76	188	6598	0.40	ng/ul	0.00
16) Chrysene-d12	20.97	240	8130	0.40	ng/ul	0.00
20) Perylene-d12	23.20	264	7583	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.72	152	2570	0.40	ng/ul	0.00
14) Fluoranthene-d10	18.80	212	8285	0.38	ng/ul	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
3) Naphthalene	10.17	128	3989	0.399	ng/ul#	94
5) 2-Methylnaphthalene	11.79	142	2532	0.401	ng/ul	97
7) Acenaphthylene	13.73	152	3252	0.394	ng/ul	98
8) Acenaphthene	14.07	153	2815	0.397	ng/ul	97
9) Fluorene	15.07	166	3277	0.403	ng/ul	93
11) Pentachlorophenol	16.44	266	219	0.337	ng/ul	95
12) Phenanthrene	16.80	178	6802	0.400	ng/ul#	89
13) Anthracene	16.90	178	5727	0.401	ng/ul#	91
15) Fluoranthene	18.83	202	8353	0.383	ng/ul	83
17) Pyrene	19.20	202	8795	0.402	ng/ul#	80
18) Benzo(a)anthracene	20.95	228	8317	0.398	ng/ul#	85
19) Chrysene	21.01	228	9612	0.399	ng/ul	97
21) Benzo(b)fluoranthene	22.53	252	7881	0.388	ng/ul	83
22) Benzo(k)fluoranthene	22.58	252	9180	0.405	ng/ul#	86
23) Benzo(a)pyrene	23.10	252	7264	0.387	ng/ul#	78
24) Indeno(1,2,3-cd)pyrene	25.50	276	9629	0.395	ng/ul#	76
25) Dibenzo(a,h)anthracene	25.52	278	8178	0.396	ng/ul#	87
26) Benzo(g,h,i)perylene	26.19	276	8197	0.395	ng/ul#	92

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

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