

Data Path : Z:\HPCHEM1\BNA E\DATA\BE031318\
 Data File : BE095754.D
 Acq On : 13 Mar 2018 17:35
 Operator : SJ/JU
 Sample : SSTDICV0.4
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SICV15

Quant Time: Mar 14 11:56:13 2018
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\SOM-EPA-SIM-BE031318.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Mar 14 11:55:57 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.42	152	1782	0.40	ng/ul	0.00
2) Naphthalene-d8	11.25	136	5055	0.40	ng/ul	0.00
6) Acenaphthene-d10	15.01	164	2799	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.71	188	6163	0.40	ng/ul	0.00
16) Chrysene-d12	21.79	240	6404	0.40	ng/ul	0.00
20) Perylene-d12	24.42	264	5746	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.80	152	3434	0.40	ng/ul	0.00
14) Fluoranthene-d10	19.68	212	7923	0.40	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	11.30	128	5754	0.408	ng/ul#	89
5) 2-Methylnaphthalene	12.87	142	3971	0.405	ng/ul	100
7) Acenaphthylene	14.73	152	5637	0.399	ng/ul	97
8) Acenaphthene	15.07	153	4335	0.407	ng/ul	92
9) Fluorene	16.03	166	4580	0.397	ng/ul#	90
11) Pentachlorophenol	17.36	266	280	0.323	ng/ul	95
12) Phenanthrene	17.75	178	7467	0.405	ng/ul#	88
13) Anthracene	17.84	178	6933	0.402	ng/ul#	91
15) Fluoranthene	19.72	202	9101	0.407	ng/ul	84
17) Pyrene	20.07	202	9769	0.402	ng/ul#	78
18) Benzo(a)anthracene	21.78	228	8214	0.398	ng/ul#	85
19) Chrysene	21.84	228	8469	0.403	ng/ul	97
21) Benzo(b)fluoranthene	23.61	252	8035	0.404	ng/ul	84
22) Benzo(k)fluoranthene	23.66	252	7691	0.406	ng/ul#	86
23) Benzo(a)pyrene	24.30	252	7381	0.406	ng/ul#	78
24) Indeno(1,2,3-cd)pyrene	27.23	276	8131	0.404	ng/ul#	77
25) Dibenzo(a,h)anthracene	27.25	278	6447	0.400	ng/ul#	88
26) Benzo(g,h,i)perylene	28.11	276	7035	0.406	ng/ul#	94

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

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