

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE072417\
 Data File : BE093535.D
 Acq On : 24 Jul 2017 13:11
 Operator : SJ/JU
 Sample : SSTDICV0.4
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SICV35

Quant Time: Jul 24 14:28:57 2017
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SOM-EPA-SIM-BE072417.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jul 24 13:14:05 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.93	152	2859	0.40	ng/ul	0.00
2) Naphthalene-d8	10.71	136	14484	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.54	164	9486	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.26	188	25709	0.40	ng/ul	0.00
16) Chrysene-d12	21.42	240	29530	0.40	ng/ul	0.00
20) Perylene-d12	23.92	264	26145	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.29	152	9618	0.41	ng/ul	0.00
14) Fluoranthene-d10	19.28	212	31109	0.38	ng/ul	0.00
Target Compounds					Qvalue	
3) Naphthalene	10.77	128	14300	0.406	ng/ul#	92
5) 2-Methylnaphthalene	12.37	142	10504	0.409	ng/ul	99
7) Acenaphthylene	14.25	152	15732	0.396	ng/ul	97
8) Acenaphthene	14.60	153	12410	0.398	ng/ul	98
9) Fluorene	15.57	166	15613	0.396	ng/ul#	92
11) Pentachlorophenol	16.93	266	1427	0.336	ng/ul	97
12) Phenanthrene	17.29	178	27702	0.396	ng/ul#	85
13) Anthracene	17.40	178	26024	0.382	ng/ul#	92
15) Fluoranthene	19.31	202	33301	0.380	ng/ul	84
17) Pyrene	19.67	202	34436	0.400	ng/ul#	83
18) Benzo(a)anthracene	21.40	228	33405	0.390	ng/ul#	86
19) Chrysene	21.45	228	32835	0.400	ng/ul	98
21) Benzo(b)fluoranthene	23.15	252	32572	0.403	ng/ul	86
22) Benzo(k)fluoranthene	23.20	252	31592	0.396	ng/ul#	88
23) Benzo(a)pyrene	23.81	252	30794	0.400	ng/ul#	82
24) Indeno(1,2,3-cd)pyrene	26.56	276	33670	0.396	ng/ul#	83
25) Dibenzo(a,h)anthracene	26.58	278	28417	0.400	ng/ul#	91
26) Benzo(g,h,i)perylene	27.38	276	28361	0.396	ng/ul	96

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

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