

Data Path : Z:\HPCHEM1\BNA E\DATA\BE101017\
 Data File : BE094271.D
 Acq On : 11 Oct 2017 3:27
 Operator : SJ/JU
 Sample : I5533-06DL 5X
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 B-77(1-3)-092817DL

Manual Integrations
 APPROVED

Sohil
 10/11/2017 7:20:31 PM

Quant Time: Oct 11 08:21:15 2017
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\SOM-EPA-SIM-BE092017.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Oct 11 07:23:41 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.91	152	2337	0.40	ng/ul	0.00
2) Naphthalene-d8	10.69	136	11627	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.51	164	6417	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.24	188	13812	0.40	ng/ul	-0.02
16) Chrysene-d12	21.41	240	14483	0.40	ng/ul	0.00
20) Perylene-d12	23.92	264	14397	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.28	152	694	0.04	ng/ul	0.00
14) Fluoranthene-d10	19.27	212	1793	0.04	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.74	128	3915	0.137	ng/ul#	94
5) 2-Methylnaphthalene	12.35	142	1324	0.072	ng/ul	98
7) Acenaphthylene	14.24	152	2758	0.110	ng/ul	99
8) Acenaphthene	14.57	153	4300	0.209	ng/ul	98
9) Fluorene	15.56	166	4809	0.199	ng/ul	93
12) Phenanthrene	17.29	178	100960	2.797	ng/ul#	92
13) Anthracene	17.38	178	24669	0.703	ng/ul#	92
15) Fluoranthene	19.30	202	187665	3.957	ng/ul	84
17) Pyrene	19.66	202	170043	4.220	ng/ul#	80
18) Benzo(a)anthracene	21.39	228	89617	2.332	ng/ul#	86
19) Chrysene	21.44	228	81607	1.839	ng/ul	99
21) Benzo(b)fluoranthene	23.15	252	101602m	2.598	ng/ul	
22) Benzo(k)fluoranthene	23.19	252	35856m	0.793	ng/ul	
23) Benzo(a)pyrene	23.80	252	75263	1.852	ng/ul#	81
24) Indeno(1,2,3-cd)pyrene	26.57	276	45250	1.088	ng/ul#	79
25) Dibenzo(a,h)anthracene	26.57	278	11486	0.329	ng/ul#	95
26) Benzo(a,h,i)perylene	27.40	276	39734	1.108	ng/ul	98

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

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