

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE110617\
 Data File : BE094765.D
 Acq On : 6 Nov 2017 22:14
 Operator : SJ/JU
 Sample : I6096-06 5X
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 B-52(1-3)-102317

Manual Integrations
 APPROVED

Sohil
 11/7/2017 5:42:00 PM

Quant Time: Nov 07 01:50:26 2017
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SOM-EPA-SIM-BE102517.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Nov 07 01:23:46 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.90	152	1426	0.40	ng/ul	-0.02
2) Naphthalene-d8	10.69	136	8288	0.40	ng/ul	-0.01
6) Acenaphthene-d10	14.51	164	4694	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.26	188	10380	0.40	ng/ul	-0.02
16) Chrysene-d12	21.43	240	10127	0.40	ng/ul	0.00
20) Perylene-d12	23.97	264	11280	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.28	152	523	0.04	ng/ul	-0.01
14) Fluoranthene-d10	19.28	212	1204	0.04	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	10.74	128	10236	0.505	ng/ul#	93
5) 2-Methylnaphthalene	12.35	142	1891	0.133	ng/ul	97
7) Acenaphthylene	14.24	152	2295	0.108	ng/ul#	94
8) Acenaphthene	14.58	153	2339	0.153	ng/ul	98
9) Fluorene	15.57	166	2787	0.163	ng/ul	95
12) Phenanthrene	17.29	178	67174m	2.450	ng/ul	
13) Anthracene	17.39	178	16042	0.593	ng/ul	96
15) Fluoranthene	19.32	202	147942	4.667	ng/ul	82
17) Pyrene	19.68	202	138081	4.411	ng/ul#	78
18) Benzo(a)anthracene	21.41	228	90600	3.129	ng/ul#	89
19) Chrysene	21.47	228	86939	2.905	ng/ul	96
21) Benzo(b)fluoranthene	23.20	252	112429m	3.543	ng/ul	
22) Benzo(k)fluoranthene	23.23	252	36697m	1.079	ng/ul	
23) Benzo(a)pyrene	23.86	252	77624	2.447	ng/ul#	86
24) Indeno(1,2,3-cd)pyrene	26.70	276	42708	1.301	ng/ul#	72
25) Dibenzo(a,h)anthracene	26.68	278	10471	0.379	ng/ul#	61
26) Benzo(g,h,i)perylene	27.54	276	33061	1.255	ng/ul#	89

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

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