

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA E\DATA\BE101417\
 Data File : BE094373.D
 Acq On : 15 Oct 2017 2:08
 Operator : SJ/JU
 Sample : I5548-03
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 C0AB4

Manual Integrations
 APPROVED

Sohil
 10/16/2017 8:25:34 PM

Quant Time: Oct 16 05:31:26 2017
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\SOM-EPA-SIM-BE092017.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Oct 16 05:09:03 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.91	152	1616	0.40	ng/ul	0.00
2) Naphthalene-d8	10.70	136	8847	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.52	164	5145	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.26	188	11122	0.40	ng/ul	0.00
16) Chrysene-d12	21.43	240	10419	0.40	ng/ul	0.00
20) Perylene-d12	23.95	264	9684m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.28	152	2462	0.18	ng/ul	0.00
14) Fluoranthene-d10	19.28	212	5745	0.15	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.75	128	1302	0.060	ng/ul#	80
5) 2-Methylnaphthalene	12.36	142	1047	0.074	ng/ul	95
7) Acenaphthylene	14.24	152	2157	0.108	ng/ul#	91
8) Acenaphthene	14.59	153	1273	0.077	ng/ul	96
9) Fluorene	15.57	166	1798	0.093	ng/ul	88
12) Phenanthrene	17.29	178	48898m	1.682	ng/ul	
13) Anthracene	17.40	178	11871	0.420	ng/ul	96
15) Fluoranthene	19.31	202	132345	3.466	ng/ul	84
17) Pyrene	19.67	202	109841	3.789	ng/ul#	83
18) Benzo(a)anthracene	21.41	228	60159	2.176	ng/ul#	89
19) Chrysene	21.46	228	67418	2.112	ng/ul	95
21) Benzo(b)fluoranthene	23.17	252	75503	2.870	ng/ul	91
22) Benzo(k)fluoranthene	23.22	252	36203m	1.191	ng/ul	
23) Benzo(a)pyrene	23.84	252	52312m	1.914	ng/ul	
24) Indeno(1,2,3-cd)pyrene	26.63	276	40103m	1.434	ng/ul	
25) Dibenzo(a,h)anthracene	26.63	278	11042m	0.471	ng/ul	
26) Benzo(a,h,i)perylene	27.47	276	33188m	1.376	ng/ul	

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

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