

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA E\DATA\BE101417\
 Data File : BE094389.D
 Acq On : 15 Oct 2017 12:35
 Operator : SJ/JU
 Sample : I5548-03DL 5X
 Misc :
 ALS Vial : 35 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 C0AB4DL

Manual Integrations
 APPROVED

Sohil
 10/16/2017 8:26:15 PM

Quant Time: Oct 16 07:22:27 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 06:50:03 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.91	152	1966	0.40	ng/ul	0.00
2) Naphthalene-d8	10.70	136	10489	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.52	164	6378	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.26	188	14946	0.40	ng/ul	0.00
16) Chrysene-d12	21.42	240	14903	0.40	ng/ul	0.00
20) Perylene-d12	23.94	264	14350	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.28	152	633	0.04	ng/ul	0.00
14) Fluoranthene-d10	19.28	212	1657	0.03	ng/ul	0.00
Target Compounds				Ovalue		
7) Acenaphthylene	14.24	152	662	0.027	ng/ul#	75
9) Fluorene	15.57	166	515	0.021	ng/ul#	85
12) Phenanthrene	17.29	178	14104	0.361	ng/ul#	89
13) Anthracene	17.38	178	3301	0.087	ng/ul#	91
15) Fluoranthene	19.31	202	38445	0.749	ng/ul	84
17) Pyrene	19.67	202	31616	0.762	ng/ul#	84
18) Benzo(a)anthracene	21.40	228	17153	0.434	ng/ul#	89
19) Chrysene	21.46	228	19902	0.436	ng/ul	96
21) Benzo(b)fluoranthene	23.17	252	25575m	0.656	ng/ul	
22) Benzo(k)fluoranthene	23.21	252	9048m	0.201	ng/ul	
23) Benzo(a)pyrene	23.83	252	15502	0.383	ng/ul#	90
24) Indeno(1,2,3-cd)pyrene	26.61	276	11318	0.273	ng/ul#	79
25) Dibenzo(a,h)anthracene	26.61	278	3042	0.088	ng/ul#	60
26) Benzo(g,h,i)perylene	27.45	276	9562	0.268	ng/ul#	91

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

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