

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA E\DATA\BE101417\
 Data File : BE094378.D
 Acq On : 15 Oct 2017 5:13
 Operator : SJ/JU
 Sample : I5548-08
 Misc :
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 C0AB9

Manual Integrations
 APPROVED

Sohil
 10/16/2017 8:25:43 PM

Quant Time: Oct 16 05:57:35 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	1504	0.40	ng/ul	0.00
2) Naphthalene-d8	10.70	136	8098	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.52	164	5181	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.26	188	11329m	0.40	ng/ul	0.00
16) Chrysene-d12	21.42	240	10171	0.40	ng/ul	0.00
20) Perylene-d12	23.95	264	11282m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.28	152	3803	0.30	ng/ul	0.00
14) Fluoranthene-d10	19.28	212	9747	0.26	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.74	128	5462	0.275	ng/ul#	95
5) 2-Methylnaphthalene	12.35	142	3840	0.298	ng/ul	98
7) Acenaphthylene	14.24	152	3280	0.162	ng/ul#	94
8) Acenaphthene	14.58	153	41512	2.496	ng/ul	97
9) Fluorene	15.57	166	48438	2.488	ng/ul#	93
12) Phenanthrene	17.29	178	788541	26.636	ng/ul#	85
13) Anthracene	17.40	178	180375	6.265	ng/ul#	91
15) Fluoranthene	19.31	202	1094648	28.140	ng/ul	84
17) Pyrene	19.68	202	862873	30.491	ng/ul#	81
18) Benzo(a)anthracene	21.40	228	438958	16.268	ng/ul#	86
19) Chrysene	21.46	228	378117	12.135	ng/ul	99
21) Benzo(b)fluoranthene	23.18	252	411257	13.419	ng/ul	85
22) Benzo(k)fluoranthene	23.22	252	175455m	4.953	ng/ul	
23) Benzo(a)pyrene	23.84	252	311035m	9.767	ng/ul	
24) Indeno(1,2,3-cd)pyrene	26.63	276	191677m	5.884	ng/ul	
25) Dibenzo(a,h)anthracene	26.63	278	52236m	1.912	ng/ul	
26) Benzo(a,h,i)perylene	27.46	276	153062m	5.447	ng/ul	

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

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