

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE110317\
 Data File : BE094718.D
 Acq On : 3 Nov 2017 21:30
 Operator : SJ/JU
 Sample : I5902-03 5X
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 B-12(0-1)-101317

Manual Integrations
 APPROVED

Sohil
 11/4/2017 2:16:40 PM

Quant Time: Nov 04 00:29:42 2017
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SOM-EPA-SIM-BE102517.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Nov 04 01:15:30 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.90	152	1471	0.40	ng/ul	-0.01
2) Naphthalene-d8	10.69	136	8864	0.40	ng/ul	-0.01
6) Acenaphthene-d10	14.51	164	4940	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.24	188	10603	0.40	ng/ul	-0.02
16) Chrysene-d12	21.42	240	10547	0.40	ng/ul	-0.01
20) Perylene-d12	23.95	264	11631	0.40	ng/ul	-0.01
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.27	152	719	0.05	ng/ul	-0.01
14) Fluoranthene-d10	19.27	212	1605	0.05	ng/ul	-0.01
Target Compounds				Qvalue		
3) Naphthalene	10.74	128	13520	0.624	ng/ul#	92
5) 2-Methylnaphthalene	12.35	142	4406	0.291	ng/ul	97
7) Acenaphthylene	14.24	152	1822	0.082	ng/ul#	90
8) Acenaphthene	14.57	153	1777	0.111	ng/ul	97
9) Fluorene	15.56	166	1884	0.104	ng/ul#	89
12) Phenanthrene	17.29	178	38957m	1.391	ng/ul	
13) Anthracene	17.38	178	7585	0.274	ng/ul	95
15) Fluoranthene	19.30	202	74315	2.295	ng/ul	84
17) Pyrene	19.67	202	65847	2.020	ng/ul#	79
18) Benzo(a)anthracene	21.40	228	42725	1.417	ng/ul#	89
19) Chrysene	21.46	228	52103	1.672	ng/ul	97
21) Benzo(b)fluoranthene	23.18	252	65645m	2.006	ng/ul	
22) Benzo(k)fluoranthene	23.22	252	20939m	0.597	ng/ul	
23) Benzo(a)pyrene	23.84	252	40392	1.235	ng/ul#	83
24) Indeno(1,2,3-cd)pyrene	26.66	276	26365	0.779	ng/ul#	75
25) Dibenzo(a,h)anthracene	26.64	278	6467	0.227	ng/ul#	82
26) Benzo(g,h,i)perylene	27.49	276	21989	0.809	ng/ul	96

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

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