

Data Path : Z:\HPCHEM1\BNA E\DATA\BE112217\
 Data File : BE094841.D
 Acq On : 22 Nov 2017 20:42
 Operator : SJ/JU
 Sample : I6496-01
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 C0AA0

Manual Integrations
 APPROVED

Sohil
 11/27/2017 3:25:37 PM

Quant Time: Nov 23 03:00:36 2017
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\SOM-EPA-SIM-BE111417.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Nov 23 02:50:08 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.90	152	795	0.40	ng/ul	-0.04
2) Naphthalene-d8	10.67	136	4885	0.40	ng/ul	-0.04
6) Acenaphthene-d10	14.50	164	3623	0.40	ng/ul	-0.01
10) Phenanthrene-d10	17.24	188	8648m	0.40	ng/ul	-0.02
16) Chrysene-d12	21.41	240	9096	0.40	ng/ul	-0.01
20) Perylene-d12	23.95	264	7486m	0.40	ng/ul	-0.02
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.27	152	2565	0.33	ng/ul	-0.03
14) Fluoranthene-d10	19.27	212	9769	0.33	ng/ul	-0.02
Target Compounds					Ovalue	
3) Naphthalene	10.72	128	2259	0.194	ng/ul#	94
5) 2-Methylnaphthalene	12.35	142	2155	0.266	ng/ul	98
7) Acenaphthylene	14.23	152	2805	0.169	ng/ul	96
8) Acenaphthene	14.56	153	745	0.064	ng/ul#	92
9) Fluorene	15.55	166	1251	0.088	ng/ul#	92
12) Phenanthrene	17.27	178	33812	1.456	ng/ul#	90
13) Anthracene	17.36	178	8356	0.353	ng/ul	96
15) Fluoranthene	19.30	202	118930	3.902	ng/ul	82
17) Pyrene	19.66	202	112501	3.952	ng/ul#	78
18) Benzo(a)anthracene	21.40	228	70620	2.554	ng/ul#	88
19) Chrysene	21.45	228	62038	2.247	ng/ul	100
21) Benzo(b)fluoranthene	23.17	252	85842m	3.849	ng/ul	
22) Benzo(k)fluoranthene	23.21	252	30474m	1.285	ng/ul	
23) Benzo(a)pyrene	23.83	252	63698m	2.872	ng/ul	
24) Indeno(1,2,3-cd)pyrene	26.65	276	39335m	1.627	ng/ul	
25) Dibenzo(a,h)anthracene	26.63	278	10519m	0.516	ng/ul	
26) Benzo(a,h,i)perylene	27.48	276	32648m	1.612	ng/ul	

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

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