

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE112817\
 Data File : BE094882.D
 Acq On : 28 Nov 2017 14:00
 Operator : SJ/JU
 Sample : I6496-01DL 5X
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 C0AA0DL

Manual Integrations
 APPROVED

Sohil
 11/29/2017 5:14:59 PM

Quant Time: Nov 29 01:47:38 2017
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SOM-EPA-SIM-BE111417.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Nov 29 01:10:42 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.91	152	562	0.40	ng/ul	-0.04
2) Naphthalene-d8	10.70	136	3838	0.40	ng/ul	-0.01
6) Acenaphthene-d10	14.51	164	2742	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.24	188	6951	0.40	ng/ul	-0.02
16) Chrysene-d12	21.42	240	6973	0.40	ng/ul	0.00
20) Perylene-d12	23.96	264	7321	0.40	ng/ul	-0.01
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.31	152	396	0.06	ng/ul	0.00
14) Fluoranthene-d10	19.28	212	1556	0.07	ng/ul	-0.01
Target Compounds				Qvalue		
3) Naphthalene	10.75	128	376	0.041	ng/ul#	62
5) 2-Methylnaphthalene	12.39	142	350	0.055	ng/ul	92
7) Acenaphthylene	14.24	152	525	0.042	ng/ul#	71
8) Acenaphthene	14.57	153	193	0.022	ng/ul#	89
9) Fluorene	15.57	166	368	0.034	ng/ul#	86
12) Phenanthrene	17.27	178	6028	0.323	ng/ul#	87
13) Anthracene	17.38	178	1989	0.105	ng/ul	96
15) Fluoranthene	19.30	202	19723	0.805	ng/ul	83
17) Pyrene	19.66	202	18781	0.861	ng/ul#	78
18) Benzo(a)anthracene	21.41	228	12293	0.580	ng/ul#	89
19) Chrysene	21.46	228	10811	0.511	ng/ul	98
21) Benzo(b)fluoranthene	23.18	252	15032m	0.689	ng/ul	
22) Benzo(k)fluoranthene	23.22	252	5772m	0.249	ng/ul	
23) Benzo(a)pyrene	23.84	252	11103	0.512	ng/ul#	80
24) Indeno(1,2,3-cd)pyrene	26.67	276	7426	0.314	ng/ul#	71
25) Dibenzo(a,h)anthracene	26.66	278	2327	0.117	ng/ul	98
26) Benzo(g,h,i)perylene	27.51	276	6317	0.319	ng/ul	93

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

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