

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE111517\
 Data File : BE094812.D
 Acq On : 15 Nov 2017 12:35
 Operator : SJ/JU
 Sample : I6260-03DL 2X
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 X2B43DL

Manual Integrations
 APPROVED

SOHIL
 11/16/2017 6:17:12 PM

Quant Time: Nov 16 11:40:36 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 16 00:58:39 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.95	152	629	0.40	ng/ul	0.01
2) Naphthalene-d8	10.75	136	3472	0.40	ng/ul	0.04
6) Acenaphthene-d10	14.55	164	2284	0.40	ng/ul	0.02
10) Phenanthrene-d10	17.27	188	6193	0.40	ng/ul	0.00
16) Chrysene-d12	21.45	240	7687	0.40	ng/ul	0.00
20) Perylene-d12	23.98	264	7537	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.37	152	945	0.17	ng/ul	0.06
14) Fluoranthene-d10	19.30	212	3647	0.17	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	10.80	128	2874	0.348	ng/ul#	87
5) 2-Methylnaphthalene	12.44	142	1292	0.225	ng/ul	98
9) Fluorene	15.61	166	2733	0.306	ng/ul	94
11) Pentachlorophenol	17.03	266	568m	1.739	ng/ul	
13) Anthracene	17.41	178	7276	0.429	ng/ul#	91
15) Fluoranthene	19.34	202	8065	0.370	ng/ul	84
17) Pyrene	19.70	202	9787	0.407	ng/ul#	81
19) Chrysene	21.49	228	7540	0.323	ng/ul	99
21) Benzo(b)fluoranthene	23.20	252	6329	0.282	ng/ul	92
23) Benzo(a)pyrene	23.88	252	8063	0.361	ng/ul#	84
25) Dibenzo(a,h)anthracene	26.70	278	15544	0.758	ng/ul	91

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

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