

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE072417\
 Data File : BE093550.D
 Acq On : 24 Jul 2017 22:23
 Operator : SJ/JU
 Sample : I4213-04
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 C0B28

Manual Integrations
 APPROVED

Sohil
 7/25/2017 5:33:13 PM

Quant Time: Jul 25 04:41:46 2017
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SOM-EPA-SIM-BE072417.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jul 24 18:07:15 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.93	152	3339	0.40	ng/ul	0.00
2) Naphthalene-d8	10.71	136	17542	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.53	164	11731	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.26	188	30133m	0.40	ng/ul	0.00
16) Chrysene-d12	21.41	240	35098	0.40	ng/ul	0.00
20) Perylene-d12	23.92	264	30386m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.29	152	7163	0.25	ng/ul	0.00
14) Fluoranthene-d10	19.28	212	25714	0.27	ng/ul	0.00
Target Compounds					Qvalue	
3) Naphthalene	10.76	128	1638	0.038	ng/ul#	91
7) Acenaphthylene	14.25	152	1531	0.031	ng/ul#	90
12) Phenanthrene	17.29	178	3041	0.037	ng/ul#	88
13) Anthracene	17.40	178	2683	0.034	ng/ul	93
15) Fluoranthene	19.31	202	28484	0.278	ng/ul	84
17) Pyrene	19.67	202	26861	0.262	ng/ul#	84
18) Benzo(a)anthracene	21.39	228	14130	0.139	ng/ul#	85
19) Chrysene	21.45	228	11687	0.120	ng/ul	98
21) Benzo(b)fluoranthene	23.15	252	15218m	0.162	ng/ul	
22) Benzo(k)fluoranthene	23.19	252	5991m	0.065	ng/ul	
23) Benzo(a)pyrene	23.81	252	10667m	0.119	ng/ul	
24) Indeno(1,2,3-cd)pyrene	26.55	276	6435m	0.065	ng/ul	
26) Benzo(g,h,i)perylene	27.37	276	5352m	0.064	ng/ul	

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

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