

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE081017\
 Data File : BE093758.D
 Acq On : 10 Aug 2017 22:41
 Operator : SJ/JU
 Sample : I4529-07
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 C0AA1

Manual Integrations
 APPROVED

sohil
 8/14/2017 7:03:50 PM

Quant Time: Aug 11 18:35:00 2017
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SOM-EPA-SIM-BE072417.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Aug 11 06:07:04 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.92	152	3019	0.40	ng/ul	0.00
2) Naphthalene-d8	10.71	136	15774	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.52	164	10084	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.26	188	21973m	0.40	ng/ul	0.00
16) Chrysene-d12	21.41	240	23417	0.40	ng/ul	0.00
20) Perylene-d12	23.91	264	19222m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.29	152	8037	0.31	ng/ul	0.00
14) Fluoranthene-d10	19.27	212	23563	0.34	ng/ul	0.00
Target Compounds						Qvalue
7) Acenaphthylene	14.24	152	4225	0.100	ng/ul	97
8) Acenaphthene	14.58	153	3024	0.091	ng/ul	96
9) Fluorene	15.57	166	5199	0.124	ng/ul	92
12) Phenanthrene	17.29	178	81386	1.360	ng/ul#	89
13) Anthracene	17.38	178	14078	0.242	ng/ul#	91
15) Fluoranthene	19.30	202	200913	2.685	ng/ul	84
17) Pyrene	19.66	202	183501	2.685	ng/ul#	84
18) Benzo(a)anthracene	21.39	228	82975	1.222	ng/ul#	86
19) Chrysene	21.44	228	85884	1.319	ng/ul	99
21) Benzo(b)fluoranthene	23.14	252	110282	1.855	ng/ul	86
22) Benzo(k)fluoranthene	23.19	252	36755m	0.627	ng/ul	
23) Benzo(a)pyrene	23.80	252	80030m	1.415	ng/ul	
24) Indeno(1,2,3-cd)pyrene	26.54	276	47805m	0.766	ng/ul	
25) Dibenzo(a,h)anthracene	26.56	278	11723m	0.225	ng/ul	
26) Benzo(g,h,i)perylene	27.36	276	38002m	0.722	ng/ul	

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

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