

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE081017\
 Data File : BE093753.D
 Acq On : 10 Aug 2017 19:00
 Operator : SJ/JU
 Sample : I4455-03
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 C0AG2

Manual Integrations
 APPROVED

sohil
 8/14/2017 7:03:09 PM

Quant Time: Aug 11 18:21:31 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	3110	0.40	ng/ul	0.00
2) Naphthalene-d8	10.71	136	16485	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.52	164	10198	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.26	188	22140	0.40	ng/ul	0.00
16) Chrysene-d12	21.41	240	23061	0.40	ng/ul	0.00
20) Perylene-d12	23.91	264	19422m	0.40	ng/ul	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
4) 2-Methylnaphthalene-d10	12.29	152	4837	0.18	ng/ul	0.00
14) Fluoranthene-d10	19.27	212	13430	0.19	ng/ul	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
7) Acenaphthylene	14.24	152	1426	0.033	ng/ul#	91
11) Pentachlorophenol	16.91	266	369	0.101	ng/ul	90
12) Phenanthrene	17.29	178	2852	0.047	ng/ul#	91
15) Fluoranthene	19.30	202	11072	0.147	ng/ul	84
17) Pyrene	19.66	202	12122	0.180	ng/ul#	83
18) Benzo(a)anthracene	21.39	228	7574	0.113	ng/ul#	86
19) Chrysene	21.45	228	7819	0.122	ng/ul	98
21) Benzo(b)fluoranthene	23.14	252	8540	0.142	ng/ul	92
22) Benzo(k)fluoranthene	23.19	252	5072m	0.086	ng/ul	
23) Benzo(a)pyrene	23.80	252	7644m	0.134	ng/ul	
24) Indeno(1,2,3-cd)pyrene	26.54	276	5242m	0.083	ng/ul	
25) Dibenzo(a,h)anthracene	26.57	278	2534m	0.048	ng/ul	
26) Benzo(g,h,i)perylene	27.37	276	4702m	0.088	ng/ul	

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

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