

Data Path : Z:\HPCHEM1\BNA E\DATA\BE101017\
 Data File : BE094264.D
 Acq On : 10 Oct 2017 23:08
 Operator : SJ/JU
 Sample : I5522-15DL 5X
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 C0AC0DL

Manual Integrations
 APPROVED

Sohil
 10/11/2017 7:20:14 PM

Quant Time: Oct 11 07:58:47 2017
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\SOM-EPA-SIM-BE092017.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Oct 11 07:23:41 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.91	152	1940	0.40	ng/ul	0.00
2) Naphthalene-d8	10.69	136	9926	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.51	164	5552	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.24	188	11976	0.40	ng/ul	-0.02
16) Chrysene-d12	21.41	240	12323	0.40	ng/ul	0.00
20) Perylene-d12	23.92	264	12357	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.28	152	1004	0.07	ng/ul	0.00
14) Fluoranthene-d10	19.27	212	2344	0.06	ng/ul	0.00
Target Compounds					Ovalue	
3) Naphthalene	10.74	128	1566	0.064	ng/ul#	91
5) 2-Methylnaphthalene	12.35	142	782	0.050	ng/ul	96
7) Acenaphthylene	14.24	152	1494	0.069	ng/ul#	92
8) Acenaphthene	14.57	153	2665	0.150	ng/ul	98
9) Fluorene	15.56	166	2724	0.131	ng/ul	93
12) Phenanthrene	17.29	178	58023	1.854	ng/ul#	93
13) Anthracene	17.38	178	14160	0.465	ng/ul#	92
15) Fluoranthene	19.30	202	134960	3.282	ng/ul	84
17) Pyrene	19.66	202	123168	3.592	ng/ul#	83
18) Benzo(a)anthracene	21.39	228	69608	2.129	ng/ul#	87
19) Chrysene	21.45	228	64673	1.713	ng/ul	99
21) Benzo(b)fluoranthene	23.15	252	90320m	2.691	ng/ul	
22) Benzo(k)fluoranthene	23.19	252	27523m	0.709	ng/ul	
23) Benzo(a)pyrene	23.80	252	62179	1.783	ng/ul#	81
24) Indeno(1,2,3-cd)pyrene	26.56	276	39945	1.119	ng/ul#	79
25) Dibenzo(a,h)anthracene	26.58	278	10111	0.338	ng/ul#	95
26) Benzo(a,h,i)perylene	27.39	276	32461	1.055	ng/ul	99

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

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