

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE081117\
 Data File : BE093780.D
 Acq On : 11 Aug 2017 21:27
 Operator : SJ/JU
 Sample : I4504-10DL 2X
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 C0AG4DL

Manual Integrations
 APPROVED

sohil
 8/14/2017 7:01:08 PM

Quant Time: Aug 12 02:59:19 2017
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SOM-EPA-SIM-BE072417.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Aug 12 02:01:53 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.92	152	2690	0.40	ng/ul	0.00
2) Naphthalene-d8	10.70	136	13942	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.52	164	8152	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.26	188	18171	0.40	ng/ul	0.00
16) Chrysene-d12	21.41	240	19177	0.40	ng/ul	0.00
20) Perylene-d12	23.91	264	18819	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.29	152	1198	0.05	ng/ul	0.00
14) Fluoranthene-d10	19.27	212	2988	0.05	ng/ul	0.00
Target Compounds				Qvalue		
8) Acenaphthene	14.58	153	753	0.028	ng/ul	97
9) Fluorene	15.57	166	2265	0.067	ng/ul	93
12) Phenanthrene	17.29	178	28643	0.579	ng/ul#	91
13) Anthracene	17.38	178	7706	0.160	ng/ul#	92
15) Fluoranthene	19.30	202	40890	0.661	ng/ul	84
17) Pyrene	19.66	202	32505	0.581	ng/ul#	85
18) Benzo(a)anthracene	21.39	228	16519	0.297	ng/ul#	87
19) Chrysene	21.44	228	15492	0.290	ng/ul	98
21) Benzo(b)fluoranthene	23.14	252	16242m	0.279	ng/ul	
22) Benzo(k)fluoranthene	23.19	252	6151m	0.107	ng/ul	
23) Benzo(a)pyrene	23.79	252	13870	0.250	ng/ul#	90
24) Indeno(1,2,3-cd)pyrene	26.54	276	6927	0.113	ng/ul#	87
25) Dibenzo(a,h)anthracene	26.56	278	3179	0.062	ng/ul#	76
26) Benzo(g,h,i)perylene	27.37	276	4663	0.091	ng/ul	93

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

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