

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE072517\
 Data File : BE093584.D
 Acq On : 25 Jul 2017 19:16
 Operator : SJ/JU
 Sample : I4213-03
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 C0B27

Manual Integrations
 APPROVED

Sohil
 7/26/2017 5:55:59 PM

Quant Time: Jul 26 03:48:05 2017
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SOM-EPA-SIM-BE072417.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jul 25 06:10:24 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.92	152	2192	0.40	ng/ul	0.00
2) Naphthalene-d8	10.71	136	11679	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.53	164	7893	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.26	188	19560m	0.40	ng/ul	0.00
16) Chrysene-d12	21.41	240	22202	0.40	ng/ul	0.00
20) Perylene-d12	23.91	264	19241m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.29	152	5155	0.27	ng/ul	0.00
14) Fluoranthene-d10	19.28	212	18223	0.30	ng/ul	0.00
Target Compounds				Qvalue		
12) Phenanthrene	17.29	178	2279	0.043	ng/ul#	89
15) Fluoranthene	19.30	202	9279	0.139	ng/ul	85
17) Pyrene	19.67	202	11278	0.174	ng/ul#	83
18) Benzo(a)anthracene	21.39	228	4800	0.075	ng/ul#	87
19) Chrysene	21.45	228	6355	0.103	ng/ul	96
21) Benzo(b)fluoranthene	23.15	252	8913	0.150	ng/ul	91
22) Benzo(k)fluoranthene	23.19	252	3156m	0.054	ng/ul	
23) Benzo(a)pyrene	23.80	252	5659m	0.100	ng/ul	
24) Indeno(1,2,3-cd)pyrene	26.55	276	3621m	0.058	ng/ul	
26) Benzo(g,h,i)perylene	27.36	276	3412m	0.065	ng/ul	

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

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