

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE102617\
 Data File : BE094593.D
 Acq On : 27 Oct 2017 13:21
 Operator : SJ/JU
 Sample : I5719-11 5X
 Misc :
 ALS Vial : 36 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 B-24(0.5-1.5)-100517

Manual Integrations
 APPROVED

Sohil
 10/27/2017 5:36:10 PM

Quant Time: Oct 27 14:50:56 2017
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SOM-EPA-SIM-BE102517.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Oct 27 04:05:19 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.91	152	1654	0.40	ng/ul	0.00
2) Naphthalene-d8	10.70	136	9516	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.52	164	5406	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.26	188	11791	0.40	ng/ul	0.00
16) Chrysene-d12	21.43	240	11573	0.40	ng/ul	0.00
20) Perylene-d12	23.97	264	11226	0.40	ng/ul	0.01
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.28	152	463	0.03	ng/ul	0.00
14) Fluoranthene-d10	19.29	212	1109	0.03	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	10.74	128	5802	0.249	ng/ul#	95
5) 2-Methylnaphthalene	12.35	142	2521	0.155	ng/ul	96
7) Acenaphthylene	14.24	152	3418	0.140	ng/ul	95
8) Acenaphthene	14.58	153	6997	0.398	ng/ul	96
9) Fluorene	15.57	166	4859	0.246	ng/ul	96
12) Phenanthrene	17.29	178	109197m	3.506	ng/ul	
13) Anthracene	17.39	178	23203	0.755	ng/ul	94
15) Fluoranthene	19.32	202	290496	8.068	ng/ul	84
17) Pyrene	19.68	202	275012	7.687	ng/ul#	80
18) Benzo(a)anthracene	21.41	228	160338	4.846	ng/ul#	88
19) Chrysene	21.47	228	162557	4.754	ng/ul	98
21) Benzo(b)fluoranthene	23.19	252	245007m	7.759	ng/ul	
22) Benzo(k)fluoranthene	23.24	252	71078m	2.100	ng/ul	
23) Benzo(a)pyrene	23.86	252	152569	4.833	ng/ul#	84
24) Indeno(1,2,3-cd)pyrene	26.68	276	101552	3.109	ng/ul#	74
25) Dibenzo(a,h)anthracene	26.68	278	24160	0.878	ng/ul#	84
26) Benzo(g,h,i)perylene	27.53	276	84393	3.218	ng/ul	96

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

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