

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE072517\
 Data File : BE093591.D
 Acq On : 25 Jul 2017 23:35
 Operator : SJ/JU
 Sample : I4209-04
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 C0B07

Manual Integrations
 APPROVED

Sohil
 7/26/2017 5:56:11 PM

Quant Time: Jul 26 05:24:15 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	2713	0.40	ng/ul	0.00
2) Naphthalene-d8	10.71	136	14472	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.53	164	9606	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.26	188	22158m	0.40	ng/ul	0.00
16) Chrysene-d12	21.41	240	23716	0.40	ng/ul	0.00
20) Perylene-d12	23.91	264	20435m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.29	152	7644	0.33	ng/ul	0.00
14) Fluoranthene-d10	19.28	212	23240	0.33	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	10.76	128	2563	0.073	ng/ul#	96
12) Phenanthrene	17.29	178	2106	0.035	ng/ul#	92
15) Fluoranthene	19.30	202	4264	0.057	ng/ul	85
17) Pyrene	19.67	202	4800	0.069	ng/ul#	88
18) Benzo(a)anthracene	21.39	228	3775	0.055	ng/ul#	88
19) Chrysene	21.45	228	4078	0.062	ng/ul	94
21) Benzo(b)fluoranthene	23.14	252	6482	0.103	ng/ul	93
22) Benzo(k)fluoranthene	23.19	252	2381m	0.038	ng/ul	
23) Benzo(a)pyrene	23.80	252	5254m	0.087	ng/ul	
24) Indeno(1,2,3-cd)pyrene	26.55	276	3436m	0.052	ng/ul	
26) Benzo(g,h,i)perylene	27.36	276	3415m	0.061	ng/ul	

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

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