

Data Path : Z:\HPCHEM1\BNA F\DATA\BF100917\
 Data File : BF099284.D
 Acq On : 10 Oct 2017 1:28
 Operator : SJ/JU
 Sample : I5715-07 25X
 Misc :
 ALS Vial : 27 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PL-02-100517-C

Manual Integrations
 APPROVED

Sohil
 10/10/2017 5:01:23 PM

Quant Time: Oct 10 05:14:36 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF092317.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Oct 05 17:20:48 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.86	152	102147	20.00	ng	0.00
21) Naphthalene-d8	8.15	136	420048	20.00	ng	0.00
38) Acenaphthene-d10	9.90	164	183002	20.00	ng	0.00
63) Phenanthrene-d10	11.39	188	279413	20.00	ng	0.00
75) Chrysene-d12	14.03	240	191739	20.00	ng	0.00
86) Perylene-d12	15.50	264	197523	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.48	112	26487	4.13	ng	0.00
7) Phenol-d6	6.49	99	33426	4.22	ng	-0.01
23) Nitrobenzene-d5	7.42	82	17802	2.72	ng	-0.01
41) 2,4,6-Tribromophenol	10.69	330	4826	3.22	ng	-0.01
44) 2-Fluorobiphenyl	9.22	172	44060	4.02	ng	-0.01
78) Terphenyl-d14	12.96	244	27965	3.32	ng	-0.01
Target Compounds						Qvalue
51) Acenaphthene	9.93	154	23475	2.050	ng	97
54) Dibenzofuran	10.10	168	34759	2.280	ng	96
57) Fluorene	10.45	166	50021	4.082	ng	96
70) Phenanthrene	11.41	178	261805	17.505	ng	99
71) Anthracene	11.46	178	101678	6.644	ng	96
74) Fluoranthene	12.60	202	251174	16.585	ng	98
77) Pyrene	12.83	202	197254	13.491	ng	98
80) Benzo(a)anthracene	14.02	228	112274	9.670	ng	99
82) Chrysene	14.05	228	84895	7.680	ng	98
85) Indeno(1,2,3-cd)pyrene	16.98	276	37969	4.294	ng	100
87) Benzo(b)fluoranthene	15.07	252	113170m	9.319	ng	
88) Benzo(k)fluoranthene	15.09	252	41451m	3.456	ng	
89) Benzo(a)pyrene	15.44	252	81665	7.326	ng	98
91) Benzo(a,h,i)perylene	17.43	276	31049	3.521	ng	98

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

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