

Data Path : Z:\HPCHEM1\BNA F\DATA\BF101817\
 Data File : BF099655.D
 Acq On : 19 Oct 2017 9:39
 Operator : SJ/JU
 Sample : I5935-06
 Misc :
 ALS Vial : 36 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 CL-02-101617-B

Manual Integrations
 APPROVED

Sohil
 10/19/2017 6:15:55 PM

Quant Time: Oct 19 17:36:42 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF092317.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Oct 18 15:18:49 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.83	152	88306	20.00	ng	0.00
21) Naphthalene-d8	8.11	136	314668	20.00	ng	0.00
38) Acenaphthene-d10	9.86	164	111406	20.00	ng	0.00
63) Phenanthrene-d10	11.35	188	179087	20.00	ng	0.00
75) Chrysene-d12	13.99	240	187342	20.00	ng	0.00
86) Perylene-d12	15.46	264	147052	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.47	112	732966	132.13	ng	0.02
7) Phenol-d6	6.47	99	776572	113.48	ng	0.00
23) Nitrobenzene-d5	7.40	82	513767	104.71	ng	0.00
41) 2,4,6-Tribromophenol	10.66	330	136880	150.15	ng	0.00
44) 2-Fluorobiphenyl	9.18	172	856951	128.41	ng	0.00
78) Terphenyl-d14	12.93	244	691080	83.89	ng	0.00
Target Compounds						
31) Naphthalene	8.13	128	226349	15.316	ng	98
37) 2-Methylnaphthalene	8.82	142	53756	5.595	ng	96
51) Acenaphthene	9.89	154	28668	4.112	ng	97
54) Dibenzofuran	10.07	168	44133	4.755	ng	97
57) Fluorene	10.41	166	23066	3.092	ng	99
70) Phenanthrene	11.37	178	41145	4.292	ng	98
72) Carbazole	11.59	167	26004	2.707	ng	95

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

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