

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF121417\
 Data File : BF101407.D
 Acq On : 14 Dec 2017 21:37
 Operator : SJ/JU
 Sample : I6841-01 100X
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 WC-01-01

Manual Integrations
 APPROVED

Sohil
 12/15/2017 6:44:25 PM

Quant Time: Dec 15 01:08:55 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF121417.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Dec 14 15:42:32 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.82	152	174615	20.00	ng	0.00
21) Naphthalene-d8	8.10	136	718693	20.00	ng	0.00
38) Acenaphthene-d10	9.85	164	322377	20.00	ng	0.00
63) Phenanthrene-d10	11.34	188	605685	20.00	ng	0.00
75) Chrysene-d12	13.99	240	411292	20.00	ng	-0.01
86) Perylene-d12	15.46	264	328070	20.00	ng	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
5) 2-Fluorophenol	0.00	112	0	0.00	ng	
7) Phenol-d6	6.55	99	1110	0.08	ng	0.09
23) Nitrobenzene-d5	7.42	82	600	0.05	ng	0.04
41) 2,4,6-Tribromophenol	10.66	330	125	0.04	ng	0.00
44) 2-Fluorobiphenyl	9.17	172	1951	0.09	ng	0.00
78) Terphenyl-d14	12.93	244	2716	0.16	ng	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
31) Naphthalene	8.12	128	779114	21.533	ng	100
37) 2-Methylnaphthalene	8.81	142	157834	6.696	ng	99
51) Acenaphthene	9.89	154	61062	2.941	ng	97
54) Dibenzofuran	10.06	168	82618	3.132	ng	98
57) Fluorene	10.40	166	136756	6.128	ng	98
70) Phenanthrene	11.37	178	564892	16.771	ng	98
71) Anthracene	11.42	178	144987	4.214	ng	98
74) Fluoranthene	12.56	202	360771	10.204	ng	96
77) Pyrene	12.79	202	344791	11.170	ng	99
80) Benzo(a)anthracene	13.98	228	125383	4.617	ng	99
82) Chrysene	14.02	228	105407	4.111	ng	97
87) Benzo(b)fluoranthene	15.04	252	77851m	3.893	ng	
89) Benzo(a)pyrene	15.40	252	73490	3.987	ng	99
91) Benzo(a,h,i)perylene	17.39	276	39080	2.494	ng	97

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

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