

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF091317\
 Data File : BF098498.D
 Acq On : 13 Sep 2017 19:17
 Operator : SJ/JU
 Sample : I5182-01
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 G-3

Manual Integrations
 APPROVED

Sohil
 9/14/2017 5:52:39 PM

Quant Time: Sep 14 06:47:39 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF091117.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Sep 12 02:14:50 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.66	152	174411	20.00	ng	-0.01
21) Naphthalene-d8	7.94	136	712565	20.00	ng	-0.01
38) Acenaphthene-d10	9.69	164	301678	20.00	ng	-0.02
63) Phenanthrene-d10	11.17	188	482354	20.00	ng	-0.02
75) Chrysene-d12	13.80	240	322176	20.00	ng	-0.01
86) Perylene-d12	15.20	264	328291	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.26	112	1102224	93.44	ng	-0.01
7) Phenol-d6	6.31	99	1383064	92.34	ng	-0.01
23) Nitrobenzene-d5	7.23	82	867187	67.85	ng	-0.01
41) 2,4,6-Tribromophenol	10.48	330	216886	107.72	ng	-0.01
44) 2-Fluorobiphenyl	9.02	172	1345244	75.58	ng	-0.02
78) Terphenyl-d14	12.75	244	820503	54.97	ng	-0.02
Target Compounds						
10) Phenol	6.32	94	54564	3.136	ng	Qvalue 88
49) Dimethylphthalate	9.41	163	185692	7.952	ng	100
70) Phenanthrene	11.19	178	603304	23.593	ng	98
71) Anthracene	11.25	178	159606	6.079	ng	97
74) Fluoranthene	12.38	202	904870	35.349	ng	99
77) Pyrene	12.61	202	836195	32.901	ng	98
80) Benzo(a)anthracene	13.79	228	428841	22.704	ng	98
82) Chrysene	13.83	228	372768	19.524	ng	96
85) Indeno(1,2,3-cd)pyrene	16.54	276	174896	13.056	ng	92
87) Benzo(b)fluoranthene	14.81	252	510560m	24.734	ng	
88) Benzo(k)fluoranthene	14.83	252	157933m	8.196	ng	
89) Benzo(a)pyrene	15.14	252	355220	19.316	ng	97
90) Dibenzo(a,h)anthracene	16.54	278	44946	3.360	ng	# 74
91) Benzo(a,h,i)perylene	16.94	276	157639	11.716	ng	95

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

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