

Data Path : Z:\HPCHEM1\BNA F\DATA\BF100317\
 Data File : BF099053.D
 Acq On : 4 Oct 2017 1:53
 Operator : SJ/JU
 Sample : I5610-03
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 2A-(1-2)-COMP-(0-5)

Manual Integrations
 APPROVED

Sohil
 10/4/2017 7:21:31 PM

Quant Time: Oct 04 04:14:54 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF092317.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Sep 23 13:50:58 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.87	152	113939	20.00	ng	-0.05
21) Naphthalene-d8	8.16	136	418860	20.00	ng	-0.05
38) Acenaphthene-d10	9.91	164	146004	20.00	ng	-0.05
63) Phenanthrene-d10	11.40	188	217200	20.00	ng	-0.05
75) Chrysene-d12	14.03	240	228415	20.00	ng	-0.05
86) Perylene-d12	15.51	264	229780	20.00	ng	-0.06
System Monitoring Compounds						
5) 2-Fluorophenol	5.50	112	969836	135.50	ng	-0.04
7) Phenol-d6	6.50	99	1120409	126.89	ng	-0.04
23) Nitrobenzene-d5	7.44	82	645994	98.91	ng	-0.05
41) 2,4,6-Tribromophenol	10.70	330	133753	111.95	ng	-0.05
44) 2-Fluorobiphenyl	9.23	172	812098	92.86	ng	-0.05
78) Terphenyl-d14	12.98	244	568126	56.57	ng	-0.05
Target Compounds						
10) Phenol	6.52	94	27662	2.801	ng	94
15) Benzyl Alcohol	7.03	79	14146	2.184	ng	# 61
49) Dimethylphthalate	9.62	163	194397	17.641	ng	98
70) Phenanthrene	11.42	178	37412	3.218	ng	97
74) Fluoranthene	12.62	202	70004	5.946	ng	95
77) Pyrene	12.84	202	65968	3.787	ng	98
80) Benzo(a)anthracene	14.02	228	36715	2.655	ng	98
82) Chrysene	14.06	228	33626	2.554	ng	98
87) Benzo(b)fluoranthene	15.07	252	42197m	2.987	ng	
89) Benzo(a)pyrene	15.44	252	33151	2.556	ng	# 93

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

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