

Data Path : Z:\HPCHEM1\BNA F\DATA\BF100317\
 Data File : BF099054.D
 Acq On : 4 Oct 2017 2:20
 Operator : SJ/JU
 Sample : I5610-04
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

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

Manual Integrations
 APPROVED

Sohil
 10/4/2017 7:21:32 PM

Quant Time: Oct 04 04:22:14 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	131910	20.00	ng	-0.05
21) Naphthalene-d8	8.16	136	477158	20.00	ng	-0.05
38) Acenaphthene-d10	9.91	164	165203	20.00	ng	-0.05
63) Phenanthrene-d10	11.40	188	248076	20.00	ng	-0.05
75) Chrysene-d12	14.03	240	266756	20.00	ng	-0.05
86) Perylene-d12	15.51	264	269623	20.00	ng	-0.06
System Monitoring Compounds						
5) 2-Fluorophenol	5.50	112	1081059	130.46	ng	-0.04
7) Phenol-d6	6.50	99	1252570	122.54	ng	-0.04
23) Nitrobenzene-d5	7.44	82	717574	96.44	ng	-0.05
41) 2,4,6-Tribromophenol	10.70	330	152409	112.74	ng	-0.05
44) 2-Fluorobiphenyl	9.23	172	928795	93.86	ng	-0.05
78) Terphenyl-d14	12.98	244	679960	57.97	ng	-0.05
Target Compounds						
10) Phenol	6.52	94	33875	2.963	ng	99
49) Dimethylphthalate	9.62	163	209359	16.791	ng	99
70) Phenanthrene	11.42	178	52734	3.971	ng	99
74) Fluoranthene	12.62	202	83677	6.223	ng	96
77) Pyrene	12.84	202	81003	3.982	ng	99
80) Benzo(a)anthracene	14.03	228	46564	2.883	ng	92
82) Chrysene	14.06	228	44371	2.885	ng	96
87) Benzo(b)fluoranthene	15.07	252	60549m	3.652	ng	
89) Benzo(a)pyrene	15.44	252	45618	2.998	ng	# 96

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

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