

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF091217\
 Data File : BF098470.D
 Acq On : 13 Sep 2017 00:29
 Operator : SJ/JU
 Sample : I5140-17
 Misc :
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 RL-13-0-16

Manual Integrations
 APPROVED

Sohil
 9/13/2017 5:00:24 PM

Quant Time: Sep 13 02:42:57 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	181688	20.00	ng	0.00
21) Naphthalene-d8	7.95	136	716982	20.00	ng	0.00
38) Acenaphthene-d10	9.69	164	275379	20.00	ng	-0.01
63) Phenanthrene-d10	11.17	188	409323	20.00	ng	-0.01
75) Chrysene-d12	13.81	240	342633	20.00	ng	0.00
86) Perylene-d12	15.21	264	293362	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.27	112	1047011	85.20	ng	0.00
7) Phenol-d6	6.32	99	1267265	81.22	ng	0.00
23) Nitrobenzene-d5	7.23	82	743059	57.78	ng	-0.01
41) 2,4,6-Tribromophenol	10.49	330	135994	73.99	ng	0.00
44) 2-Fluorobiphenyl	9.02	172	955406	58.80	ng	-0.01
78) Terphenyl-d14	12.76	244	592280	37.31	ng	-0.01
Target Compounds						
10) Phenol	6.32	94	47064	2.597	ng	Qvalue # 73
49) Dimethylphthalate	9.42	163	152233	7.142	ng	98
70) Phenanthrene	11.20	178	128304	5.913	ng	98
74) Fluoranthene	12.39	202	268409	12.356	ng	97
77) Pyrene	12.61	202	228638	8.459	ng	98
80) Benzo(a)anthracene	13.80	228	125071	6.226	ng	96
82) Chrysene	13.83	228	116300	5.727	ng	96
85) Indeno(1,2,3-cd)pyrene	16.55	276	37436	2.628	ng	98
87) Benzo(b)fluoranthene	14.82	252	124864m	6.769	ng	
88) Benzo(k)fluoranthene	14.84	252	39899m	2.317	ng	
89) Benzo(a)pyrene	15.15	252	81203	4.941	ng	# 93
91) Benzo(a,h,i)perylene	16.96	276	34127	2.838	ng	# 90

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

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