

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF110117\
 Data File : BF100064.D
 Acq On : 2 Nov 2017 00:09
 Operator : SJ/JU
 Sample : I6247-01
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 S13-SP3-C

Manual Integrations
 APPROVED

Sohil
 11/2/2017 4:27:44 PM

Quant Time: Nov 02 04:56:26 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF102317.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Nov 01 18:13:08 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.77	152	80831	20.00	ng	0.00
21) Naphthalene-d8	8.06	136	313195	20.00	ng	0.00
38) Acenaphthene-d10	9.82	164	133004	20.00	ng	0.00
63) Phenanthrene-d10	11.32	188	204358	20.00	ng	0.01
75) Chrysene-d12	14.28	240	149033m	20.00	ng	0.31
86) Perylene-d12	16.12	264	142117m	20.00	ng	0.66
System Monitoring Compounds						
5) 2-Fluorophenol	5.40	112	439505	85.36	ng	0.00
7) Phenol-d6	6.42	99	526273	86.21	ng	0.00
23) Nitrobenzene-d5	7.35	82	285930	58.78	ng	0.00
41) 2,4,6-Tribromophenol	10.62	330	93352	77.02	ng	0.00
44) 2-Fluorobiphenyl	9.14	172	422713	49.03	ng	0.00
78) Terphenyl-d14	13.00	244	264011m	38.71	ng	0.10
Target Compounds						
10) Phenol	6.44	94	25736	3.840	ng	89
49) Dimethylphthalate	9.53	163	126309	11.992	ng	99
70) Phenanthrene	11.34	178	94153	8.164	ng	98
74) Fluoranthene	12.60	202	129391	10.796	ng	97
77) Pyrene	12.85	202	114302	10.086	ng	99
80) Benzo(a)anthracene	14.26	228	70183m	7.429	ng	
82) Chrysene	14.31	228	74288m	8.364	ng	
87) Benzo(b)fluoranthene	15.64	252	110890m	12.357	ng	
88) Benzo(k)fluoranthene	15.67	252	42222m	4.989	ng	
89) Benzo(a)pyrene	15.99	252	54759m	6.793	ng	
91) Benzo(g,h,i)perylene	17.70	276	43209m	6.166	ng	

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

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