

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF112316\
 Data File : BF091296.D
 Acq On : 24 Nov 2016 1:05
 Operator : UM/SJ
 Sample : H5740-12
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-05(0-0.16)

Manual Integrations
 APPROVED

sohil
 11/28/2016 4:54:09 PM

Quant Time: Nov 24 05:41:50 2016
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF112316.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Nov 23 13:22:53 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.88	152	204911	20.00	ng	0.00
21) Naphthalene-d8	8.17	136	856756	20.00	ng	0.00
38) Acenaphthene-d10	9.92	164	402464	20.00	ng	0.00
63) Phenanthrene-d10	11.40	188	564694	20.00	ng	0.00
75) Chrysene-d12	14.04	240	485217	20.00	ng	0.00
86) Perylene-d12	15.48	264	413750	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.49	112	1597405	125.25	ng	0.02
7) Phenol-d6	6.51	99	2325199	141.90	ng	0.01
23) Nitrobenzene-d5	7.45	82	1451963	97.56	ng	0.00
41) 2,4,6-Tribromophenol	10.70	330	466319	114.90	ng	0.00
44) 2-Fluorobiphenyl	9.24	172	2040872	83.03	ng	0.00
78) Terphenyl-d14	12.99	244	1234196	58.09	ng	0.00
Target Compounds						
10) Phenol	6.52	94	80955	4.43	ng	# 23
49) Dimethylphthalate	9.63	163	118967	4.73	ng	# 97
70) Phenanthrene	11.42	178	391115	13.35	ng	99
71) Anthracene	11.47	178	107053	3.57	ng	98
74) Fluoranthene	12.61	202	711472	23.88	ng	99
77) Pyrene	12.84	202	618850	17.26	ng	99
80) Benzo(a)anthracene	14.03	228	343205	12.46	ng	97
82) Chrysene	14.07	228	300309m	11.54	ng	
85) Indeno(1,2,3-cd)pyrene	16.89	276	148818m	6.64	ng	
87) Benzo(b)fluoranthene	15.06	252	445810m	17.08	ng	
88) Benzo(k)fluoranthene	15.08	252	81924m	3.84	ng	
89) Benzo(a)pyrene	15.41	252	259941m	11.71	ng	
91) Benzo(g,h,i)perylene	17.31	276	127300	6.26	ng	97

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

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