

Data Path : Z:\HPCHEM1\BNA F\DATA\BF110716\
 Data File : BF091061.D
 Acq On : 7 Nov 2016 17:38
 Operator : UM/SJ
 Sample : H5539-10
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 V001-BF017-01

Manual Integrations
 APPROVED

sohil
 11/8/2016 4:32:02 PM

Quant Time: Nov 08 00:33:45 2016
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF110316.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Nov 08 00:06:22 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.67	152	187729	20.00	ng	0.00
21) Naphthalene-d8	7.95	136	723214	20.00	ng	-0.01
38) Acenaphthene-d10	9.71	164	408322	20.00	ng	0.00
63) Phenanthrene-d10	11.19	188	594273	20.00	ng	-0.01
75) Chrysene-d12	13.83	240	408858	20.00	ng	0.00
86) Perylene-d12	15.21	264	367865	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.25	112	1163402	102.35	ng	0.00
7) Phenol-d6	6.33	99	1702212	110.33	ng	0.00
23) Nitrobenzene-d5	7.24	82	1086932	81.99	ng	0.00
41) 2,4,6-Tribromophenol	10.50	330	337833	91.52	ng	-0.01
44) 2-Fluorobiphenyl	9.04	172	1420578	59.90	ng	-0.01
78) Terphenyl-d14	12.77	244	1237700	75.54	ng	-0.01
Target Compounds						
10) Phenol	6.34	94	93640	5.48	ng	Qvalue # 53
49) Dimethylphthalate	9.44	163	97602	3.65	ng	98
70) Phenanthrene	11.21	178	277035	8.77	ng	99
74) Fluoranthene	12.40	202	440277	13.44	ng	96
77) Pyrene	12.63	202	369134	13.79	ng	98
80) Benzo(a)anthracene	13.81	228	155685m	6.71	ng	
82) Chrysene	13.85	228	145868m	6.37	ng	
85) Indeno(1,2,3-cd)pyrene	16.50	276	81295	4.28	ng	# 90
87) Benzo(b)fluoranthene	14.82	252	166195m	6.66	ng	
88) Benzo(k)fluoranthene	14.83	252	76374m	3.49	ng	
89) Benzo(a)pyrene	15.15	252	124870	5.75	ng	# 97
91) Benzo(a,h,i)perylene	16.88	276	75442	4.44	ng	98

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

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