

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF021121\
 Data File : BF123187.D
 Acq On : 11 Feb 2021 15:58
 Operator : JU/CG
 Sample : M1380-01
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PM-WC-C-001

Manual Integrations
 APPROVED

mohammad
 2/12/2021 11:23:57 AM

Quant Time: Feb 11 16:39:09 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF012721.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Feb 11 14:10:38 2021
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.881	152	124847	20.00	ng	0.00
21) Naphthalene-d8	8.163	136	496858	20.00	ng	0.00
39) Acenaphthene-d10	9.922	164	268902	20.00	ng	0.00
64) Phenanthrene-d10	11.416	188	487709	20.00	ng	0.00
76) Chrysene-d12	14.063	240	388973	20.00	ng	-0.01
86) Perylene-d12	15.551	264	348049	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.493	112	165543	20.45	ng	0.00
7) Phenol-d6	6.498	99	612164	55.80	ng	-0.01
23) Nitrobenzene-d5	7.439	82	517354	52.40	ng	0.00
42) 2,4,6-Tribromophenol	10.716	330	8364	2.87	ng	0.00
45) 2-Fluorobiphenyl	9.245	172	980702	54.22	ng	0.00
79) Terphenyl-d14	13.004	244	1103485	55.74	ng	0.00
Target Compounds						Qvalue
10) Phenol	6.510	94	22111	2.03	ng	93
50) Dimethylphthalate	9.633	163	51038	2.42	ng	100
71) Phenanthrene	11.439	178	195616	6.46	ng	97
75) Fluoranthene	12.633	202	193774	6.40	ng	97
78) Pyrene	12.863	202	143643	4.89	ng	96
81) Benzo(a)anthracene	14.051	228	71748	2.73	ng	97
83) Chrysene	14.086	228	64834	2.46	ng	94
88) Benzo(b)fluoranthene	15.110	252	61728m	2.46	ng	

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

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