

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF112120\
 Data File : BF122437.D
 Acq On : 21 Nov 2020 23:58
 Operator : JU/CG
 Sample : L4808-01RE
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 11931RE

Manual Integrations
 APPROVED

mohammad
 11/24/2020 9:02:29 AM

Quant Time: Nov 23 01:02:29 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF110920.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Nov 20 13:27:02 2020
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.863	152	201537	20.00	ng	0.00
21) Naphthalene-d8	8.151	136	794077	20.00	ng	# 0.00
39) Acenaphthene-d10	9.904	164	415738	20.00	ng	-0.01
64) Phenanthrene-d10	11.392	188	758113	20.00	ng	-0.01
76) Chrysene-d12	14.039	240	626337	20.00	ng	-0.01
86) Perylene-d12	15.521	264	524190	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.481	112	300569	22.90	ng	-0.01
7) Phenol-d6	6.492	99	1000414	58.26	ng	-0.02
23) Nitrobenzene-d5	7.422	82	875855	67.52	ng	-0.01
42) 2,4,6-Tribromophenol	10.692	330	31422	7.04	ng	-0.01
45) 2-Fluorobiphenyl	9.227	172	1575338	59.21	ng	0.00
79) Terphenyl-d14	12.986	244	1694329	57.31	ng	0.00
Target Compounds						
						Qvalue
10) Phenol	6.504	94	84649	5.01	ng	76
50) Dimethylphthalate	9.622	163	72621	2.46	ng	97
71) Phenanthrene	11.416	178	177726	4.13	ng	98
75) Fluoranthene	12.610	202	195059	4.34	ng	97
78) Pyrene	12.839	202	167624	3.81	ng	98
88) Benzo(b)fluoranthene	15.086	252	68808m	2.03	ng	

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

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