

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF120820\
 Data File : BF122590.D
 Acq On : 8 Dec 2020 16:10
 Operator : JU/CG
 Sample : L4970-12
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 4S1-A-D-COMP

Manual Integrations
 APPROVED

mohammad
 12/9/2020 12:26:37 PM

Quant Time: Dec 08 17:13:40 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF120520.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Dec 07 12:08:49 2020
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.840	152	156769	20.00	ng	0.00
21) Naphthalene-d8	8.122	136	625208	20.00	ng	0.00
39) Acenaphthene-d10	9.875	164	332297	20.00	ng	0.00
64) Phenanthrene-d10	11.363	188	569076	20.00	ng	0.00
76) Chrysene-d12	14.010	240	384592	20.00	ng	0.00
86) Perylene-d12	15.468	264	384752	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.457	112	944054	95.34	ng	0.00
7) Phenol-d6	6.469	99	1194519	90.96	ng	-0.01
23) Nitrobenzene-d5	7.398	82	718922	57.61	ng	-0.01
42) 2,4,6-Tribromophenol	10.669	330	340817	78.36	ng	0.00
45) 2-Fluorobiphenyl	9.198	172	1330754	54.17	ng	0.00
79) Terphenyl-d14	12.957	244	1180050	53.71	ng	0.00
Target Compounds						
10) Phenol	6.481	94	49408	3.63	ng	Qvalue 87
50) Dimethylphthalate	9.592	163	132502	4.99	ng	99
75) Fluoranthene	12.580	202	97514	2.75	ng	77
78) Pyrene	12.810	202	83527	2.81	ng	# 86
88) Benzo(b)fluoranthene	15.045	252	63995m	2.37	ng	

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

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