

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF120920\
 Data File : BF122610.D
 Acq On : 9 Dec 2020 13:34
 Operator : JU/CG
 Sample : L4969-01
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 NORTH-BRANFORD-01

Manual Integrations
 APPROVED

mohammad
 12/10/2020 2:42:46 PM

Quant Time: Dec 09 17:12:24 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	6.839	152	174237	20.00	ng	0.00
21) Naphthalene-d8	8.122	136	671844	20.00	ng	0.00
39) Acenaphthene-d10	9.875	164	349266	20.00	ng	0.00
64) Phenanthrene-d10	11.363	188	606669	20.00	ng	0.00
76) Chrysene-d12	14.004	240	454496	20.00	ng	0.00
86) Perylene-d12	15.468	264	465860	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.463	112	1050525	95.45	ng	0.00
7) Phenol-d6	6.469	99	1327192	90.93	ng	-0.01
23) Nitrobenzene-d5	7.398	82	836860	62.41	ng	-0.01
42) 2,4,6-Tribromophenol	10.663	330	288559	63.12	ng	-0.01
45) 2-Fluorobiphenyl	9.198	172	1532156	59.33	ng	0.00
79) Terphenyl-d14	12.951	244	1422778	54.80	ng	0.00
Target Compounds						Qvalue
10) Phenol	6.481	94	37547	2.48	ng	91
32) Benzoic acid	7.869	122	20758m	2.73	ng	
50) Dimethylphthalate	9.592	163	106478	3.82	ng	100
71) Phenanthrene	11.386	178	76011	2.11	ng	96
75) Fluoranthene	12.574	202	131606	3.48	ng	98
78) Pyrene	12.804	202	140038	3.99	ng	99
88) Benzo(b)fluoranthene	15.039	252	76527m	2.34	ng	

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

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