

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF121020\
 Data File : BF122650.D
 Acq On : 10 Dec 2020 17:12
 Operator : JU/CG
 Sample : L4990-02
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SIDEWALK

Manual Integrations
 APPROVED

mohammad
 12/11/2020 1:21:21 PM

Quant Time: Dec 11 00:34:41 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	158986	20.00	ng	0.00
21) Naphthalene-d8	8.122	136	617774	20.00	ng	0.00
39) Acenaphthene-d10	9.874	164	329857	20.00	ng	0.00
64) Phenanthrene-d10	11.363	188	609520	20.00	ng	0.00
76) Chrysene-d12	13.998	240	567694	20.00	ng	-0.01
86) Perylene-d12	15.462	264	490561	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.451	112	94049	9.37	ng	0.00
7) Phenol-d6	6.463	99	687136	51.59	ng	-0.02
23) Nitrobenzene-d5	7.398	82	655420	53.16	ng	-0.01
42) 2,4,6-Tribromophenol	10.674	330	424	0.10	ng	0.00
45) 2-Fluorobiphenyl	9.198	172	1271472	52.14	ng	0.00
79) Terphenyl-d14	12.945	244	1472052	45.39	ng	-0.01
Target Compounds						Qvalue
10) Phenol	6.475	94	42227	3.06	ng	91
50) Dimethylphthalate	9.592	163	72071	2.74	ng	98
71) Phenanthrene	11.386	178	213973	5.91	ng	97
75) Fluoranthene	12.574	202	244265	6.43	ng	96
78) Pyrene	12.798	202	215146	4.90	ng	99
81) Benzo(a)anthracene	13.986	228	102978	2.71	ng	98
83) Chrysene	14.021	228	100013	2.65	ng	96
88) Benzo(b)fluoranthene	15.033	252	88978m	2.59	ng	
90) Benzo(a)pyrene	15.398	252	62025	2.06	ng	95

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

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