

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF120820\
 Data File : BF122589.D
 Acq On : 8 Dec 2020 15:38
 Operator : JU/CG
 Sample : L4965-11
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 CRT014

Manual Integrations
 APPROVED

mohammad
 12/9/2020 12:26:35 PM

Quant Time: Dec 08 16:35:26 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	169044	20.00	ng	0.00
21) Naphthalene-d8	8.122	136	656789	20.00	ng	0.00
39) Acenaphthene-d10	9.874	164	356954	20.00	ng	0.00
64) Phenanthrene-d10	11.363	188	672068	20.00	ng	0.00
76) Chrysene-d12	14.004	240	521352	20.00	ng	0.00
86) Perylene-d12	15.462	264	464460	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.457	112	781950	73.23	ng	0.00
7) Phenol-d6	6.469	99	987921	69.76	ng	-0.01
23) Nitrobenzene-d5	7.398	82	602513	45.96	ng	-0.01
42) 2,4,6-Tribromophenol	10.663	330	248111	53.10	ng	-0.01
45) 2-Fluorobiphenyl	9.198	172	990237	37.52	ng	0.00
79) Terphenyl-d14	12.945	244	1047103	35.16	ng	-0.01
Target Compounds						Qvalue
10) Phenol	6.475	94	44928	3.06	ng	95
50) Dimethylphthalate	9.592	163	111890	3.92	ng	99
71) Phenanthrene	11.386	178	89392	2.24	ng	98
75) Fluoranthene	12.574	202	160327	3.83	ng	98
78) Pyrene	12.804	202	152107	3.77	ng	98
88) Benzo(b)fluoranthene	15.039	252	69862m	2.14	ng	

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

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