

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF112520\
 Data File : BF122487.D
 Acq On : 25 Nov 2020 16:25
 Operator : JU/CG
 Sample : L4882-06
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PAV-B48-112320-10-16.5

Manual Integrations
 APPROVED

mohammad
 11/27/2020 11:42:46 AM

Quant Time: Nov 25 17:06:46 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF110920.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Nov 24 14:38:56 2020
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.857	152	203848	20.00	ng	0.00
21) Naphthalene-d8	8.139	136	791275	20.00	ng	0.00
39) Acenaphthene-d10	9.898	164	413995	20.00	ng	0.00
64) Phenanthrene-d10	11.380	188	723467	20.00	ng	0.00
76) Chrysene-d12	14.021	240	486254	20.00	ng	-0.01
86) Perylene-d12	15.492	264	497367	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.492	112	1385037	104.33	ng	0.01
7) Phenol-d6	6.492	99	1577831	90.85	ng	0.00
23) Nitrobenzene-d5	7.422	82	1198582	92.73	ng	0.00
42) 2,4,6-Tribromophenol	10.686	330	575779	129.58	ng	0.00
45) 2-Fluorobiphenyl	9.222	172	2150820	81.17	ng	0.00
79) Terphenyl-d14	12.974	244	2075274	90.42	ng	0.00
Target Compounds						
10) Phenol	6.504	94	49266	2.88	ng	Qvalue 80
32) Benzoic acid	7.881	122	4421m	8.45	ng	

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

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