

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF071420\
 Data File : BF120863.D
 Acq On : 14 Jul 2020 16:18
 Operator : JU/CG
 Sample : L3282-02
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 ST-JOES-AVE-2

Manual Integrations
 APPROVED

mohammad
 7/15/2020 10:59:47 AM

Quant Time: Jul 14 16:44:28 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF070220.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jul 13 14:16:14 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.83	152	204479	20.00	ng	0.00
21) Naphthalene-d8	8.11	136	786889	20.00	ng	0.00
39) Acenaphthene-d10	9.87	164	410022	20.00	ng	0.00
64) Phenanthrene-d10	11.35	188	767091	20.00	ng	0.00
76) Chrysene-d12	13.99	240	606161	20.00	ng	-0.02
86) Perylene-d12	15.44	264	497934	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.44	112	701401	52.95	ng	0.00
7) Phenol-d6	6.45	99	1053095	62.25	ng	-0.01
23) Nitrobenzene-d5	7.39	82	714190	52.58	ng	0.00
42) 2,4,6-Tribromophenol	10.65	330	171324	41.59	ng	-0.01
45) 2-Fluorobiphenyl	9.19	172	1308641	48.29	ng	0.00
79) Terphenyl-d14	12.94	244	1293826	41.79	ng	0.00
Target Compounds						
50) Dimethylphthalate	9.58	163	161121	5.387	ng	98
71) Phenanthrene	11.37	178	85946	2.065	ng	97
75) Fluoranthene	12.56	202	201704	4.549	ng	97
78) Pyrene	12.79	202	181936	3.758	ng	97
81) Benzo(a)anthracene	13.97	228	87491	2.252	ng	98
83) Chrysene	14.01	228	88575	2.235	ng	96
88) Benzo(b)fluoranthene	15.02	252	104557m	3.194	ng	

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

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