

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF072318\
 Data File : BF107582.D
 Acq On : 23 Jul 2018 12:37
 Operator : JU/SJ
 Sample : J3820-05
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 BU-01-071918

Manual Integrations
 APPROVED

Sohil
 7/24/2018 5:25:36 PM

Quant Time: Jul 23 14:38:16 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF072018.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jul 20 16:29:27 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.97	152	349512	20.00	ng	0.00
21) Naphthalene-d8	8.25	136	1357456	20.00	ng	0.00
38) Acenaphthene-d10	10.01	164	627983	20.00	ng	0.00
63) Phenanthrene-d10	11.51	188	1225927	20.00	ng	0.00
75) Chrysene-d12	14.16	240	777644	20.00	ng	0.00
86) Perylene-d12	15.69	264	832261	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.61	112	2054865	94.53	ng	0.00
7) Phenol-d6	6.61	99	2485608	95.23	ng	0.00
23) Nitrobenzene-d5	7.54	82	1639159	81.49	ng	-0.01
41) 2,4,6-Tribromophenol	10.81	330	780683	109.32	ng	0.00
44) 2-Fluorobiphenyl	9.33	172	2651171	68.46	ng	0.00
78) Terphenyl-d14	13.10	244	2315207	76.22	ng	0.00
Target Compounds						
10) Phenol	6.62	94	82923	2.701	ng	55
49) Dimethylphthalate	9.72	163	232675	4.916	ng	99
70) Phenanthrene	11.54	178	171763	2.875	ng	97
74) Fluoranthene	12.73	202	262990	4.102	ng	96
77) Pyrene	12.96	202	298199	5.606	ng	98
80) Benzo(a)anthracene	14.15	228	96925	2.127	ng	96
82) Chrysene	14.18	228	116587	2.663	ng	98
87) Benzo(b)fluoranthene	15.24	252	112796m	2.476	ng	
89) Benzo(a)pyrene	15.62	252	94351	2.265	ng	98

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

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