

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF062118\
 Data File : BF106663.D
 Acq On : 21 Jun 2018 15:53
 Operator : JU/SJ
 Sample : J3524-01 2X
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 AREA-3(S)

Manual Integrations
 APPROVED

Sohil
 6/22/2018 10:23:02 AM

Quant Time: Jun 21 16:28:25 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF061918.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jun 19 14:32:09 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.96	152	96089	20.00	ng	0.00
21) Naphthalene-d8	8.24	136	381516	20.00	ng	0.00
38) Acenaphthene-d10	10.00	164	179036	20.00	ng	0.00
63) Phenanthrene-d10	11.50	188	300239	20.00	ng	0.00
75) Chrysene-d12	14.15	240	262887	20.00	ng	0.00
86) Perylene-d12	15.70	264	198011	20.00	ng	0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.60	112	130521	23.00	ng	0.00
7) Phenol-d6	6.60	99	171036	24.14	ng	-0.01
23) Nitrobenzene-d5	7.52	82	103057	15.01	ng	-0.01
41) 2,4,6-Tribromophenol	10.80	330	31162	15.02	ng	0.00
44) 2-Fluorobiphenyl	9.31	172	203426	5.42	ng	-0.01
78) Terphenyl-d14	13.08	244	192756	17.76	ng	0.00
Target Compounds						
70) Phenanthrene	11.52	178	86244	5.956	ng	98
74) Fluoranthene	12.71	202	107597	6.741	ng	93
77) Pyrene	12.94	202	102705	5.925	ng	99
80) Benzo(a)anthracene	14.14	228	50701	3.164	ng	97
82) Chrysene	14.18	228	48518	3.292	ng	97
87) Benzo(b)fluoranthene	15.24	252	43301m	3.520	ng	
89) Benzo(a)pyrene	15.63	252	28893	2.642	ng	# 92

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

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