

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF061318\
 Data File : BF106393.D
 Acq On : 13 Jun 2018 16:11
 Operator : JU/SJ
 Sample : J3458-01
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 EPE-105-5(0-5)

Manual Integrations
 APPROVED

Sohil
 6/14/2018 2:47:40 PM

Quant Time: Jun 14 03:12:11 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF061118.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jun 13 10:51:26 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.97	152	139299	20.00	ng	-0.01
21) Naphthalene-d8	8.25	136	527886	20.00	ng	-0.01
38) Acenaphthene-d10	10.01	164	256970	20.00	ng	-0.01
63) Phenanthrene-d10	11.50	188	508357	20.00	ng	0.00
75) Chrysene-d12	14.15	240	375554	20.00	ng	-0.01
86) Perylene-d12	15.68	264	384781	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.61	112	845172	102.69	ng	-0.02
7) Phenol-d6	6.61	99	1065187	102.44	ng	-0.02
23) Nitrobenzene-d5	7.52	82	672646	74.96	ng	-0.01
41) 2,4,6-Tribromophenol	10.80	330	308615	124.82	ng	-0.01
44) 2-Fluorobiphenyl	9.32	172	998727	63.34	ng	0.00
78) Terphenyl-d14	13.08	244	1138906	75.32	ng	0.00
Target Compounds						
49) Dimethylphthalate	9.71	163	205824	11.211	ng	# 90
57) Fluorene	10.55	166	41388	2.464	ng	# 92
70) Phenanthrene	11.52	178	106333	4.271	ng	96
74) Fluoranthene	12.72	202	267956	10.092	ng	96
77) Pyrene	12.94	202	252035	10.719	ng	99
80) Benzo(a)anthracene	14.14	228	147641	6.606	ng	98
82) Chrysene	14.17	228	118998	5.832	ng	97
85) Indeno(1,2,3-cd)pyrene	17.24	276	63141	3.877	ng	92
87) Benzo(b)fluoranthene	15.22	252	151109m	6.497	ng	
89) Benzo(a)pyrene	15.62	252	112116	5.359	ng	96
91) Benzo(g,h,i)perylene	17.71	276	58672	3.461	ng	# 93

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

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