

Data Path : U:\HPCHEM1\BNA F\DATA\BF011818\
 Data File : BF102200.D
 Acq On : 18 Jan 2018 19:13
 Operator : SJ/JU
 Sample : J1168-08 2X
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-57-0-2.0-DUP

Manual Integrations
 APPROVED

Sohil
 1/19/2018 9:47:32 AM

Quant Time: Jan 19 01:23:23 2018
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF010918.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jan 17 15:28:18 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.72	152	164026	20.00	ng	0.00
21) Naphthalene-d8	8.00	136	647923	20.00	ng	0.00
38) Acenaphthene-d10	9.76	164	277243	20.00	ng	0.00
63) Phenanthrene-d10	11.24	188	425207	20.00	ng	0.00
75) Chrysene-d12	13.87	240	281713	20.00	ng	0.00
86) Perylene-d12	15.29	264	263277	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.32	112	510956	43.99	ng	0.00
7) Phenol-d6	6.35	99	611433	44.12	ng	0.00
23) Nitrobenzene-d5	7.28	82	368715	33.44	ng	-0.01
41) 2,4,6-Tribromophenol	10.54	330	115800	47.86	ng	0.00
44) 2-Fluorobiphenyl	9.08	172	680006	38.32	ng	0.00
78) Terphenyl-d14	12.82	244	457083	33.69	ng	0.00
Target Compounds						
49) Dimethylphthalate	9.47	163	89539	4.026	ng	99
70) Phenanthrene	11.26	178	90886	3.659	ng	98
74) Fluoranthene	12.44	202	163307	6.683	ng	97
77) Pyrene	12.67	202	147755	5.934	ng	98
80) Benzo(a)anthracene	13.86	228	74063	4.121	ng	97
82) Chrysene	13.89	228	69056	3.685	ng	97
83) Bis(2-ethylhexyl)phthalate	13.86	149	30560	2.331	ng	97
85) Indeno(1,2,3-cd)pyrene	16.66	276	34488	2.528	ng	98
87) Benzo(b)fluoranthene	14.88	252	89507m	5.214	ng	
88) Benzo(k)fluoranthene	14.90	252	33146m	2.002	ng	
89) Benzo(a)pyrene	15.23	252	64822	4.196	ng	# 91
91) Benzo(a,h,i)perylene	17.08	276	36018	2.901	ng	93

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

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