

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF092718\
 Data File : BF109407.D
 Acq On : 28 Sep 2018 1:45
 Operator : JU/SJ
 Sample : J5084-21
 Misc :
 ALS Vial : 28 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 B-13(1-1.5)

Manual Integrations
 APPROVED

Sohil
 9/28/2018 2:40:34 PM

Quant Time: Sep 28 07:19:15 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF092218.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Sep 22 13:32:18 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.70	152	123424	20.00	ng	0.00
21) Naphthalene-d8	7.99	136	471038	20.00	ng	0.00
38) Acenaphthene-d10	9.74	164	221688	20.00	ng	0.00
63) Phenanthrene-d10	11.22	188	376829	20.00	ng	0.00
75) Chrysene-d12	13.84	240	267896	20.00	ng	0.00
86) Perylene-d12	15.24	264	232977	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.32	112	631521	84.28	ng	0.00
7) Phenol-d6	6.35	99	816844	85.43	ng	0.00
23) Nitrobenzene-d5	7.27	82	536382	67.22	ng	-0.01
41) 2,4,6-Tribromophenol	10.52	330	196744	90.03	ng	-0.01
44) 2-Fluorobiphenyl	9.06	172	952088	64.77	ng	-0.01
78) Terphenyl-d14	12.80	244	706396	64.71	ng	0.00
Target Compounds						
49) Dimethylphthalate	9.46	163	301666	18.709	ng	100
70) Phenanthrene	11.24	178	117666	6.254	ng	97
74) Fluoranthene	12.42	202	147745	7.348	ng	95
77) Pyrene	12.64	202	150854	7.857	ng	98
80) Benzo(a)anthracene	13.83	228	61880	3.835	ng	96
82) Chrysene	13.86	228	70280	4.394	ng	98
87) Benzo(b)fluoranthene	14.84	252	69944m	5.464	ng	
89) Benzo(a)pyrene	15.18	252	46476	4.029	ng	99
91) Benzo(g,h,i)perylene	16.99	276	24885	2.676	ng	91

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

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