

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF092518\
 Data File : BF109306.D
 Acq On : 26 Sep 2018 00:20
 Operator : JU/SJ
 Sample : J5044-07
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 GRID07-COMP-092018

Manual Integrations
 APPROVED

Sohil
 9/26/2018 3:43:01 PM

Quant Time: Sep 26 04:57:06 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	6.71	152	114580	20.00	ng	0.00
21) Naphthalene-d8	7.99	136	426518	20.00	ng	0.00
38) Acenaphthene-d10	9.74	164	181953	20.00	ng	0.00
63) Phenanthrene-d10	11.22	188	295869	20.00	ng	0.00
75) Chrysene-d12	13.84	240	266805	20.00	ng	0.00
86) Perylene-d12	15.24	264	196409	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.32	112	483414	69.49	ng	0.00
7) Phenol-d6	6.35	99	618109	69.63	ng	-0.01
23) Nitrobenzene-d5	7.27	82	378597	52.40	ng	-0.01
41) 2,4,6-Tribromophenol	10.52	330	115757	64.54	ng	-0.01
44) 2-Fluorobiphenyl	9.06	172	633609	52.52	ng	-0.01
78) Terphenyl-d14	12.80	244	512300	47.12	ng	0.00
Target Compounds						
10) Phenol	6.36	94	27199	2.706	ng	98
49) Dimethylphthalate	9.46	163	182775	13.811	ng	99
70) Phenanthrene	11.24	178	35246	2.386	ng	96
74) Fluoranthene	12.42	202	50482	3.198	ng	96
77) Pyrene	12.65	202	44941	2.350	ng	97
87) Benzo(b)fluoranthene	14.85	252	24318m	2.253	ng	

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

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