

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF092518\
 Data File : BF109299.D
 Acq On : 25 Sep 2018 21:15
 Operator : JU/SJ
 Sample : J5044-08
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 GRID08-COMP-092018

Manual Integrations
 APPROVED

Sohil
 9/26/2018 3:42:17 PM

Quant Time: Sep 26 10:23:41 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	111407	20.00	ng	0.00
21) Naphthalene-d8	7.99	136	426880	20.00	ng	0.00
38) Acenaphthene-d10	9.74	164	188239	20.00	ng	0.00
63) Phenanthrene-d10	11.22	188	325360	20.00	ng	0.00
75) Chrysene-d12	13.84	240	277139	20.00	ng	0.00
86) Perylene-d12	15.24	264	212696	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.32	112	430376	63.63	ng	0.00
7) Phenol-d6	6.35	99	554984	64.30	ng	-0.01
23) Nitrobenzene-d5	7.27	82	334092	46.20	ng	-0.01
41) 2,4,6-Tribromophenol	10.52	330	110113	59.34	ng	-0.01
44) 2-Fluorobiphenyl	9.06	172	599359	48.02	ng	-0.01
78) Terphenyl-d14	12.80	244	541259	47.93	ng	0.00
Target Compounds						
10) Phenol	6.36	94	25812	2.641	ng	97
49) Dimethylphthalate	9.46	163	176611	12.900	ng	99
74) Fluoranthene	12.42	202	46851	2.699	ng	96
77) Pyrene	12.65	202	41848	2.107	ng	98
80) Benzo(a)anthracene	13.83	228	42474	2.544	ng	98
87) Benzo(b)fluoranthene	14.85	252	37053m	3.170	ng	
89) Benzo(a)pyrene	15.19	252	27229	2.585	ng	93

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

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