

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF100118\
 Data File : BF109476.D
 Acq On : 1 Oct 2018 16:54
 Operator : JU/SJ
 Sample : J4780-03
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TR-04-092818

Manual Integrations
 APPROVED

Sohil
 10/2/2018 1:01:28 PM

Quant Time: Oct 01 18:31:07 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	119177	20.00	ng	0.00
21) Naphthalene-d8	7.99	136	445134	20.00	ng	0.00
38) Acenaphthene-d10	9.73	164	184167	20.00	ng	0.00
63) Phenanthrene-d10	11.22	188	318214	20.00	ng	0.00
75) Chrysene-d12	13.84	240	261326	20.00	ng	0.00
86) Perylene-d12	15.25	264	186697	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.33	112	615533	85.07	ng	0.01
7) Phenol-d6	6.36	99	762771	82.61	ng	0.00
23) Nitrobenzene-d5	7.27	82	475962	63.12	ng	-0.01
41) 2,4,6-Tribromophenol	10.52	330	154287	84.98	ng	-0.01
44) 2-Fluorobiphenyl	9.06	172	790013	64.70	ng	-0.01
78) Terphenyl-d14	12.80	244	635312	59.66	ng	0.00
Target Compounds						
49) Dimethylphthalate	9.46	163	79017	5.899	ng	Qvalue # 99
70) Phenanthrene	11.24	178	79469	5.002	ng	99
74) Fluoranthene	12.43	202	146633	8.636	ng	96
77) Pyrene	12.65	202	267406	14.278	ng	99
80) Benzo(a)anthracene	13.83	228	130227	8.273	ng	93
82) Chrysene	13.87	228	139835	8.963	ng	97
85) Indeno(1,2,3-cd)pyrene	16.60	276	41764	3.303	ng	93
87) Benzo(b)fluoranthene	14.85	252	100176m	9.765	ng	
88) Benzo(k)fluoranthene	14.87	252	35163m	3.612	ng	
89) Benzo(a)pyrene	15.19	252	84339	9.124	ng	97
91) Benzo(g,h,i)perylene	17.02	276	49009	6.575	ng	# 90

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

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