

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF092818\
 Data File : BF109431.D
 Acq On : 28 Sep 2018 15:54
 Operator : JU/SJ
 Sample : J4776-07 2X
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 EO-01-092718-B

Manual Integrations
 APPROVED

Sohil
 10/1/2018 9:57:01 AM

Quant Time: Sep 28 18:47:18 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	126337	20.00	ng	0.00
21) Naphthalene-d8	7.99	136	455279	20.00	ng	0.00
38) Acenaphthene-d10	9.74	164	194688	20.00	ng	0.00
63) Phenanthrene-d10	11.22	188	290197	20.00	ng	0.00
75) Chrysene-d12	13.84	240	240171	20.00	ng	0.00
86) Perylene-d12	15.25	264	191353	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.32	112	294467	38.39	ng	0.00
7) Phenol-d6	6.35	99	375151	38.33	ng	-0.01
23) Nitrobenzene-d5	7.26	82	232218	30.11	ng	-0.02
41) 2,4,6-Tribromophenol	10.52	330	51630	26.90	ng	-0.01
44) 2-Fluorobiphenyl	9.06	172	401070	31.07	ng	-0.01
78) Terphenyl-d14	12.80	244	270183	27.61	ng	0.00
Target Compounds						
49) Dimethylphthalate	9.46	163	58072	4.101	ng	Qvalue # 98
70) Phenanthrene	11.24	178	85222	5.882	ng	98
74) Fluoranthene	12.43	202	101737	6.570	ng	89
77) Pyrene	12.66	202	119491	6.942	ng	99
80) Benzo(a)anthracene	13.83	228	63048	4.358	ng	99
82) Chrysene	13.87	228	50476	3.520	ng	96
87) Benzo(b)fluoranthene	14.85	252	55060m	5.236	ng	
89) Benzo(a)pyrene	15.19	252	36446	3.847	ng	# 90
91) Benzo(g,h,i)perylene	17.00	276	16164	2.116	ng	# 94

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

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