

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF111122\
 Data File : BF131159.D
 Acq On : 11 Nov 2022 18:11
 Operator : CG\JU
 Sample : N5506-13
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TP-10

Manual Integrations
 APPROVED

Reviewed By : Christian Giraldo 11/12/2022
 Supervised By : Jagrut Upadhyay 11/12/2022

Quant Time: Nov 12 00:05:43 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF110922.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Nov 12 00:03:31 2022
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.869	152	90602	20.000	ng	0.00
21) Naphthalene-d8	8.157	136	389288	20.000	ng	0.00
39) Acenaphthene-d10	9.916	164	159531	20.000	ng	0.00
64) Phenanthrene-d10	11.404	188	295075	20.000	ng	0.00
76) Chrysene-d12	14.051	240	140594	20.000	ng	0.00
86) Perylene-d12	15.545	264	136143	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.492	112	482268	90.523	ng	0.00
7) Phenol-d6	6.504	99	667384	95.817	ng	0.00
23) Nitrobenzene-d5	7.439	82	418835	48.543	ng	0.00
42) 2,4,6-Tribromophenol	10.710	330	160921	97.604	ng	0.00
45) 2-Fluorobiphenyl	9.233	172	662525	57.974	ng	0.00
79) Terphenyl-d14	12.992	244	473442	59.042	ng	0.00
Target Compounds						Qvalue
71) Phenanthrene	11.427	178	59538	3.669	ng	98
75) Fluoranthene	12.621	202	72852	4.485	ng	99
78) Pyrene	12.851	202	80228	6.651	ng	99
81) Benzo(a)anthracene	14.039	228	31938	3.178	ng	92
83) Chrysene	14.074	228	30904	3.256	ng	96
87) Indeno(1,2,3-cd)pyrene	17.045	276	15443	5.140	ng	92
88) Benzo(b)fluoranthene	15.104	252	35220m	3.659	ng	
90) Benzo(a)pyrene	15.474	252	27999	3.628	ng	# 94
91) Dibenzo(a,h)anthracene	17.051	278	4774	4.326	ng	# 72
92) Benzo(g,h,i)perylene	17.498	276	18819	6.153	ng	96

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

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