

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF111318\
 Data File : BF110729.D
 Acq On : 13 Nov 2018 19:54
 Operator : JU/SJ
 Sample : J5925-03
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TP-1S

Manual Integrations
 APPROVED

Sohil
 11/14/2018 1:52:22 PM

Quant Time: Nov 14 05:52:42 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF110818.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Nov 13 12:43:10 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.95	152	85015	20.00	ng	0.00
21) Naphthalene-d8	8.23	136	338552	20.00	ng	0.00
39) Acenaphthene-d10	9.99	164	145642	20.00	ng	0.00
64) Phenanthrene-d10	11.48	188	220980	20.00	ng	0.00
76) Chrysene-d12	14.13	240	158192	20.00	ng	-0.01
87) Perylene-d12	15.66	264	147350	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.57	112	590857	112.89	ng	0.00
7) Phenol-d6	6.57	99	747211	111.64	ng	0.00
23) Nitrobenzene-d5	7.51	82	469742	79.91	ng	0.00
42) 2,4,6-Tribromophenol	10.78	330	144731	98.62	ng	0.00
45) 2-Fluorobiphenyl	9.30	172	763105	74.83	ng	0.00
79) Terphenyl-d14	13.06	244	491259	58.16	ng	-0.01
Target Compounds						
10) Phenol	6.58	94	32814	4.228	ng	Qvalue # 74
50) Dimethylphthalate	9.69	163	128589	11.833	ng	99
71) Phenanthrene	11.50	178	24914	2.104	ng	99
75) Fluoranthene	12.70	202	36414	3.068	ng	99
78) Pyrene	12.93	202	38681	3.176	ng	99
81) Benzo(a)anthracene	14.12	228	21688	2.294	ng	95
83) Chrysene	14.16	228	20561	2.283	ng	96
88) Benzo(b)fluoranthene	15.20	252	21010m	2.383	ng	
90) Benzo(a)pyrene	15.59	252	16471	2.057	ng	# 94

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

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