

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF121418\
 Data File : BF111424.D
 Acq On : 14 Dec 2018 17:44
 Operator : JU/SJ
 Sample : J6319-14
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 068_DUP-2

Manual Integrations
 APPROVED

Sohil
 12/17/2018 2:36:32 PM

Quant Time: Dec 14 18:13:22 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF121418.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Dec 14 13:26:17 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.79	152	173377	20.00	ng	0.00
21) Naphthalene-d8	8.07	136	695072	20.00	ng	0.00
39) Acenaphthene-d10	9.83	164	307316	20.00	ng	0.00
64) Phenanthrene-d10	11.31	188	436803	20.00	ng	0.00
76) Chrysene-d12	13.95	240	330573	20.00	ng	0.00
87) Perylene-d12	15.39	264	271298	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.42	112	920954	88.39	ng	0.00
7) Phenol-d6	6.44	99	1179251	85.09	ng	0.00
23) Nitrobenzene-d5	7.35	82	717443	61.47	ng	0.00
42) 2,4,6-Tribromophenol	10.62	330	197395	71.70	ng	0.00
45) 2-Fluorobiphenyl	9.15	172	1181237	59.37	ng	0.00
79) Terphenyl-d14	12.90	244	708585	50.33	ng	0.00
Target Compounds						
10) Phenol	6.45	94	84145	5.450	ng	92
50) Dimethylphthalate	9.55	163	110251	4.981	ng	99
71) Phenanthrene	11.33	178	384125	17.066	ng	94
72) Anthracene	11.39	178	75876	3.296	ng	94
75) Fluoranthene	12.53	202	578559	26.275	ng	96
78) Pyrene	12.75	202	645873	27.712	ng	96
81) Benzo(a)anthracene	13.94	228	313357	17.433	ng	97
83) Chrysene	13.97	228	297303	15.701	ng	97
86) Indeno(1,2,3-cd)pyrene	16.80	276	116823	8.548	ng	92
88) Benzo(b)fluoranthene	14.98	252	350952m	22.188	ng	
89) Benzo(k)fluoranthene	15.00	252	85514m	5.658	ng	
90) Benzo(a)pyrene	15.33	252	242049	16.970	ng	97
91) Dibenzo(a,h)anthracene	16.82	278	29247	2.599	ng	# 88
92) Benzo(a,h,i)perylene	17.23	276	116640	10.561	ng	93

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

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