

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF121918\
 Data File : BF111602.D
 Acq On : 19 Dec 2018 20:48
 Operator : JU/SJ
 Sample : J6403-57
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TP-8-A

Manual Integrations
 APPROVED

Sohil
 12/20/2018 12:42:08 PM

Quant Time: Dec 20 03:56:23 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF121918.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Dec 19 17:39:39 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.89	152	192179	20.00	ng	0.00
21) Naphthalene-d8	8.17	136	725622	20.00	ng	0.00
39) Acenaphthene-d10	9.93	164	369242	20.00	ng	0.00
64) Phenanthrene-d10	11.41	188	687169	20.00	ng	0.00
76) Chrysene-d12	14.04	240	520472	20.00	ng	-0.01
87) Perylene-d12	15.52	264	528821	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.52	112	920682	81.66	ng	0.00
7) Phenol-d6	6.52	99	1196790	83.41	ng	0.00
23) Nitrobenzene-d5	7.45	82	718344	57.62	ng	0.00
42) 2,4,6-Tribromophenol	10.71	330	347907	79.74	ng	-0.01
45) 2-Fluorobiphenyl	9.25	172	1298041	51.24	ng	0.00
79) Terphenyl-d14	12.99	244	1320981	48.13	ng	0.00
Target Compounds						
10) Phenol	6.53	94	64588	3.989	ng	93
50) Dimethylphthalate	9.64	163	134514	4.982	ng	98
75) Fluoranthene	12.62	202	151102	4.094	ng	94
78) Pyrene	12.84	202	143388	3.901	ng	95
81) Benzo(a)anthracene	14.03	228	72031	2.425	ng	98
83) Chrysene	14.06	228	66671	2.284	ng	96
88) Benzo(b)fluoranthene	15.08	252	86436m	2.797	ng	
90) Benzo(a)pyrene	15.44	252	64224m	2.287	ng	

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

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