

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF010719\
 Data File : BF111879.D
 Acq On : 7 Jan 2019 16:32
 Operator : JU/SJ
 Sample : K1041-01
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 OS-1-5-COMP

Manual Integrations
 APPROVED

Sohil
 1/8/2019 11:20:46 AM

Quant Time: Jan 07 17:11:30 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF010719.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jan 07 14:33:51 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.89	152	180297	20.00	ng	0.00
21) Naphthalene-d8	8.17	136	695889	20.00	ng	0.00
39) Acenaphthene-d10	9.93	164	331743	20.00	ng	0.00
64) Phenanthrene-d10	11.41	188	609417	20.00	ng	0.00
76) Chrysene-d12	14.05	240	399102	20.00	ng	0.00
87) Perylene-d12	15.54	264	517170	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.53	112	1116668	105.96	ng	0.00
7) Phenol-d6	6.53	99	1364264	100.44	ng	0.00
23) Nitrobenzene-d5	7.45	82	906258	69.60	ng	-0.01
42) 2,4,6-Tribromophenol	10.72	330	402434	93.26	ng	0.00
45) 2-Fluorobiphenyl	9.25	172	1608820	70.72	ng	0.00
79) Terphenyl-d14	13.00	244	1411724	67.14	ng	0.00
Target Compounds						
10) Phenol	6.54	94	69660	4.719	ng	84
50) Dimethylphthalate	9.63	163	146284	5.870	ng	98
71) Phenanthrene	11.43	178	99001	3.038	ng	99
75) Fluoranthene	12.62	202	164319	4.913	ng	99
78) Pyrene	12.85	202	141832	5.165	ng	96
81) Benzo(a)anthracene	14.04	228	67350	2.932	ng	98
83) Chrysene	14.07	228	60439	2.760	ng	98
86) Indeno(1,2,3-cd)pyrene	17.03	276	40422	2.532	ng	98
88) Benzo(b)fluoranthene	15.11	252	100291m	3.242	ng	
90) Benzo(a)pyrene	15.47	252	73655	2.708	ng	# 98

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

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