

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF053019\
 Data File : BF114685.D
 Acq On : 31 May 2019 2:32
 Operator : HP/JU
 Sample : K2642-07 2X
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 OK-02-052919

Manual Integrations
 APPROVED

mohammad
 6/3/2019 8:47:59 AM

Quant Time: May 31 07:56:39 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF052919.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed May 29 19:08:51 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.70	152	394604	20.00	ng	0.00
21) Naphthalene-d8	7.97	136	1410667	20.00	ng	0.00
39) Acenaphthene-d10	9.72	164	617652	20.00	ng	0.00
64) Phenanthrene-d10	11.20	188	1110350	20.00	ng	0.00
76) Chrysene-d12	13.82	240	680778	20.00	ng	0.00
87) Perylene-d12	15.22	264	582601	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.30	112	1152324	51.27	ng	0.00
7) Phenol-d6	6.33	99	1393766	51.46	ng	0.00
23) Nitrobenzene-d5	7.25	82	773700	34.22	ng	0.00
42) 2,4,6-Tribromophenol	10.50	330	282031	47.92	ng	0.00
45) 2-Fluorobiphenyl	9.05	172	1427666	33.78	ng	0.00
79) Terphenyl-d14	12.78	244	1317997	36.40	ng	0.00
Target Compounds						
50) Dimethylphthalate	9.45	163	219264	5.095	ng	98
71) Phenanthrene	11.22	178	266805	4.710	ng	96
75) Fluoranthene	12.40	202	716092	12.210	ng	96
78) Pyrene	12.63	202	720911	12.518	ng	98
81) Benzo(a)anthracene	13.81	228	344987	6.594	ng	95
83) Chrysene	13.84	228	325356	6.509	ng	95
86) Indeno(1,2,3-cd)pyrene	16.54	276	163246m	3.261	ng	
88) Benzo(b)fluoranthene	14.82	252	364018m	9.920	ng	
89) Benzo(k)fluoranthene	14.85	252	134848m	4.228	ng	
90) Benzo(a)pyrene	15.16	252	264399	8.202	ng	# 87
92) Benzo(g,h,i)perylene	16.94	276	161102	5.827	ng	97

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

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