

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF030519\
 Data File : BF112860.D
 Acq On : 5 Mar 2019 18:53
 Operator : JU/SJ
 Sample : K1715-11
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TP-5-A

Manual Integrations
 APPROVED

Sohil
 3/6/2019 12:24:27 PM

Quant Time: Mar 06 03:49:08 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF022719.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Mar 01 15:40:43 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.74	152	156046	20.00	ng	0.00
21) Naphthalene-d8	8.02	136	608204	20.00	ng	0.00
39) Acenaphthene-d10	9.77	164	301585	20.00	ng	0.00
64) Phenanthrene-d10	11.25	188	633193	20.00	ng	0.00
76) Chrysene-d12	13.87	240	449896	20.00	ng	0.00
87) Perylene-d12	15.28	264	434786	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.35	112	1060831	105.75	ng	0.00
7) Phenol-d6	6.37	99	1413913	112.20	ng	0.00
23) Nitrobenzene-d5	7.30	82	868243	85.42	ng	0.00
42) 2,4,6-Tribromophenol	10.56	330	468274	146.26	ng	0.00
45) 2-Fluorobiphenyl	9.09	172	1498965	81.44	ng	0.00
79) Terphenyl-d14	12.83	244	1615594	69.61	ng	0.00
Target Compounds						
50) Dimethylphthalate	9.49	163	157354	7.076	ng	99
71) Phenanthrene	11.27	178	74978	2.380	ng	95
75) Fluoranthene	12.46	202	158668	4.701	ng	96
78) Pyrene	12.68	202	137710	3.631	ng	97
81) Benzo(a)anthracene	13.86	228	69624	2.233	ng	93
83) Chrysene	13.90	228	87867m	2.877	ng	
86) Indeno(1,2,3-cd)pyrene	16.64	276	52104	2.001	ng	# 91
88) Benzo(b)fluoranthene	14.88	252	128746m	4.578	ng	
90) Benzo(a)pyrene	15.22	252	64305	2.585	ng	96
92) Benzo(g,h,i)perylene	17.04	276	44825	2.103	ng	# 95

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

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