

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF060119\
 Data File : BF114747.D
 Acq On : 1 Jun 2019 10:53
 Operator : HP/JU
 Sample : K3097-09
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 0429-HR-9(HACKENSACK)

Manual Integrations
 APPROVED

mohammad
 6/3/2019 8:51:09 AM

Quant Time: Jun 03 02:36:39 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	6.69	152	388666	20.00	ng	0.00
21) Naphthalene-d8	7.97	136	1475404	20.00	ng	0.00
39) Acenaphthene-d10	9.72	164	722000	20.00	ng	0.00
64) Phenanthrene-d10	11.19	188	1427692	20.00	ng	0.00
76) Chrysene-d12	13.82	240	1023706	20.00	ng	0.00
87) Perylene-d12	15.20	264	946180	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.31	112	2510225	113.39	ng	0.02
7) Phenol-d6	6.33	99	3099669	116.20	ng	0.00
23) Nitrobenzene-d5	7.25	82	1939450	82.01	ng	0.00
42) 2,4,6-Tribromophenol	10.50	330	796096	115.72	ng	0.00
45) 2-Fluorobiphenyl	9.05	172	2968383	75.43	ng	0.00
79) Terphenyl-d14	12.78	244	3214254	59.03	ng	0.00
Target Compounds						
10) Phenol	6.34	94	131708	4.087	ng	82
50) Dimethylphthalate	9.44	163	503560	10.009	ng	98
75) Fluoranthene	12.40	202	439300	5.825	ng	93
78) Pyrene	12.62	202	355354	4.104	ng	95
88) Benzo(b)fluoranthene	14.81	252	139358m	2.338	ng	

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

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