

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP101819\
 Data File : BP000883.D
 Acq On : 18 Oct 2019 18:14
 Operator : JU
 Sample : K5394-04 2X
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 T-15-E1-(0-20)

Manual Integrations
 APPROVED

mohammad
 10/22/2019 1:39:23 AM

Quant Time: Oct 19 05:04:13 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP100119.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Oct 18 16:40:44 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.74	152	157675	20.00	ng	0.00
21) Naphthalene-d8	8.02	136	575902	20.00	ng	0.00
39) Acenaphthene-d10	9.79	164	310880	20.00	ng	0.00
64) Phenanthrene-d10	11.29	188	569186	20.00	ng	0.00
76) Chrysene-d12	13.96	240	503502	20.00	ng	0.00
87) Perylene-d12	15.45	264	569963	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.36	112	259137	26.42	ng	0.00
7) Phenol-d6	6.38	99	523216	44.26	ng	0.00
23) Nitrobenzene-d5	7.30	82	295537	31.34	ng	0.00
42) 2,4,6-Tribromophenol	10.58	330	24358	7.35	ng	0.00
45) 2-Fluorobiphenyl	9.10	172	716581	37.29	ng	0.00
79) Terphenyl-d14	12.89	244	831612	33.44	ng	0.00
Target Compounds						
50) Dimethylphthalate	9.50	163	94047	4.354	ng	99
71) Phenanthrene	11.31	178	256652	8.978	ng	99
72) Anthracene	11.36	178	84277	2.973	ng	99
75) Fluoranthene	12.52	202	687640	21.418	ng	98
78) Pyrene	12.75	202	605521	17.437	ng	98
81) Benzo(a)anthracene	13.96	228	288974	9.406	ng	98
83) Chrysene	13.99	228	256825	8.518	ng	97
86) Indeno(1,2,3-cd)pyrene	16.92	276	85147	2.261	ng	96
88) Benzo(b)fluoranthene	15.02	252	254720m	7.877	ng	
89) Benzo(k)fluoranthene	15.05	252	88460m	2.873	ng	
90) Benzo(a)pyrene	15.38	252	178950	5.937	ng	# 89
92) Benzo(a,h,i)perylene	17.37	276	80810	2.720	ng	# 90

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

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