

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP091819\
 Data File : BP000310.D
 Acq On : 18 Sep 2019 20:27
 Operator : HP/JU
 Sample : K4668-08 2X
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SU-01-091719

Manual Integrations
 APPROVED

Jagrut
 9/19/2019 11:07:17 AM

Quant Time: Sep 19 04:50:47 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP091619.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Sep 16 18:59:05 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.79	152	317087	20.00	ng	0.00
21) Naphthalene-d8	8.08	136	1150892	20.00	ng	0.00
39) Acenaphthene-d10	9.83	164	624108	20.00	ng	0.00
64) Phenanthrene-d10	11.33	188	1073826	20.00	ng	0.00
76) Chrysene-d12	14.00	240	812088	20.00	ng	0.00
87) Perylene-d12	15.49	264	880041	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.42	112	838681	45.66	ng	0.01
7) Phenol-d6	6.43	99	1071251	47.58	ng	0.00
23) Nitrobenzene-d5	7.35	82	556313	31.19	ng	0.00
42) 2,4,6-Tribromophenol	10.63	330	326018	52.67	ng	0.00
45) 2-Fluorobiphenyl	9.15	172	1340419	38.66	ng	0.00
79) Terphenyl-d14	12.93	244	1375877	35.54	ng	0.00
Target Compounds						
50) Dimethylphthalate	9.55	163	172825	4.064	ng	98
71) Phenanthrene	11.35	178	148776	2.765	ng	99
75) Fluoranthene	12.56	202	1004060	17.376	ng	96
78) Pyrene	12.79	202	1098310	18.077	ng	99
81) Benzo(a)anthracene	13.99	228	878160	17.119	ng	95
83) Chrysene	14.02	228	724652	14.388	ng	97
86) Indeno(1,2,3-cd)pyrene	16.97	276	518589m	8.447	ng	
88) Benzo(b)fluoranthene	15.05	252	1179894m	22.474	ng	
89) Benzo(k)fluoranthene	15.07	252	417545m	9.118	ng	
90) Benzo(a)pyrene	15.42	252	867506	18.160	ng	99
91) Dibenzo(a,h)anthracene	16.99	278	132486	2.975	ng	# 85
92) Benzo(a,h,i)perylene	17.43	276	447007	9.764	ng	97

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

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