

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP010620\
 Data File : BP001543.D
 Acq On : 06 Jan 2020 22:42
 Operator : JU
 Sample : K6463-05
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 GP-5C-(5-6)

Manual Integrations
 APPROVED

mohammad
 1/7/2020 12:31:46 PM

Quant Time: Jan 07 03:57:06 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP010320.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jan 06 16:38:29 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.68	152	56841	20.00	ng	0.00
21) Naphthalene-d8	10.46	136	207211	20.00	ng	0.00
39) Acenaphthene-d10	14.33	164	125432	20.00	ng	0.00
64) Phenanthrene-d10	17.09	188	267155	20.00	ng	0.00
76) Chrysene-d12	21.19	240	226955	20.00	ng	0.00
87) Perylene-d12	23.44	264	238709	20.00	ng	0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.32	112	283196	83.70	ng	0.00
7) Phenol-d6	6.89	99	413489	90.73	ng	0.00
23) Nitrobenzene-d5	8.83	82	287890	68.62	ng	0.00
42) 2,4,6-Tribromophenol	15.83	330	152156	107.18	ng	0.00
45) 2-Fluorobiphenyl	12.95	172	502536	58.91	ng	0.00
79) Terphenyl-d14	19.67	244	708927	62.25	ng	0.00
Target Compounds						
10) Phenol	6.91	94	11437	2.386	ng	93
50) Dimethylphthalate	13.79	163	41515	4.234	ng	# 88
71) Phenanthrene	17.13	178	34108	2.354	ng	# 83
75) Fluoranthene	19.11	202	123067	7.016	ng	98
78) Pyrene	19.46	202	217447	13.203	ng	98
81) Benzo(a)anthracene	21.17	228	39182	2.460	ng	# 62
83) Chrysene	21.22	228	48085	3.149	ng	# 78
88) Benzo(b)fluoranthene	22.77	252	55032m	3.723	ng	
90) Benzo(a)pyrene	23.34	252	31193	2.255	ng	# 57
92) Benzo(g,h,i)perylene	26.42	276	39522	2.882	ng	# 92

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

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