

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP010820\
 Data File : BP001570.D
 Acq On : 09 Jan 2020 01:21
 Operator : JU
 Sample : L1051-05
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SB-4F-(22-24)

Manual Integrations
 APPROVED

mohammad
 1/9/2020 12:12:10 PM

Quant Time: Jan 09 05:43:31 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP010820.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jan 08 15:02:19 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.66	152	32940	20.00	ng	0.00
21) Naphthalene-d8	10.43	136	138404	20.00	ng	0.00
39) Acenaphthene-d10	14.30	164	101654	20.00	ng	0.00
64) Phenanthrene-d10	17.06	188	248217	20.00	ng	-0.01
76) Chrysene-d12	21.16	240	232831	20.00	ng	-0.01
87) Perylene-d12	23.42	264	244669	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.29	112	257405	153.08	ng	0.00
7) Phenol-d6	6.86	99	396568	147.56	ng	0.00
23) Nitrobenzene-d5	8.80	82	286245	90.37	ng	0.00
42) 2,4,6-Tribromophenol	15.80	330	206565	133.49	ng	0.00
45) 2-Fluorobiphenyl	12.92	172	609092	88.10	ng	0.00
79) Terphenyl-d14	19.65	244	1002297	78.74	ng	0.00
Target Compounds						
10) Phenol	6.88	94	9528	3.427	ng	89
50) Dimethylphthalate	13.77	163	24240	2.895	ng	98
75) Fluoranthene	19.08	202	69252	3.899	ng	97
78) Pyrene	19.43	202	56828	3.323	ng	97
81) Benzo(a)anthracene	21.15	228	33474	2.064	ng	97
88) Benzo(b)fluoranthene	22.74	252	42404m	2.750	ng	

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

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